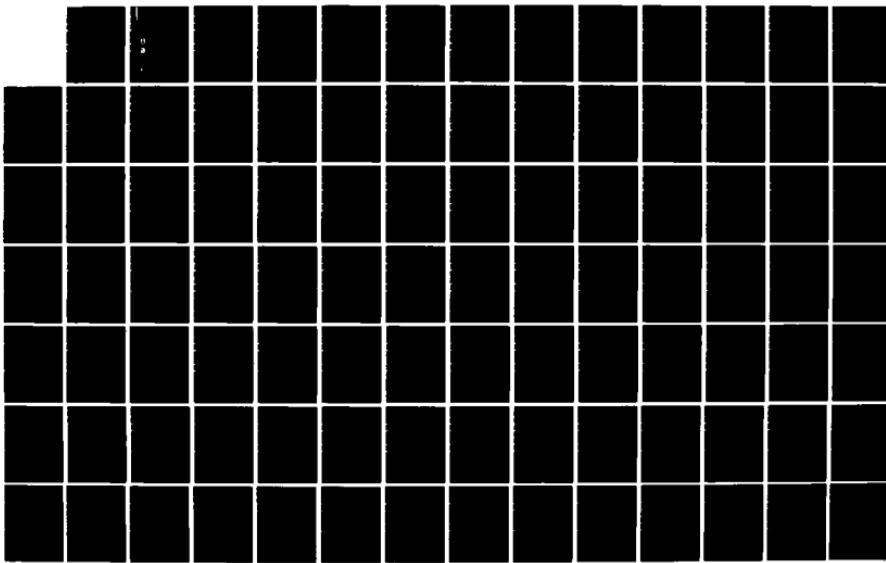
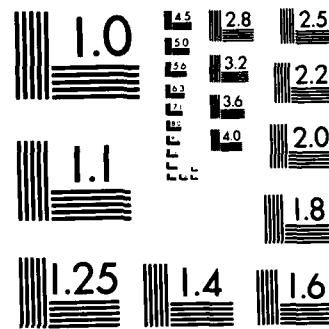


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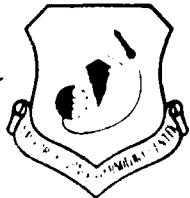
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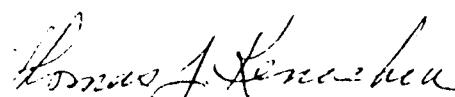
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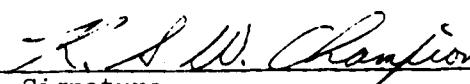
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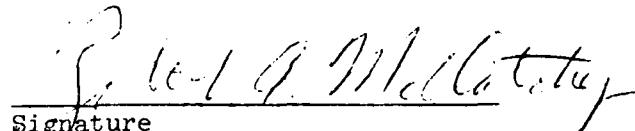
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Model of the Mesosphere and Lower Thermosphere

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A Computer Code for a One-Dimensional Dynamic Model of the Mesosphere and Lower Thermosphere

I. INTRODUCTION

T.J. Keneshea, one of the authors, pioneered studies of the numerical solution of sets of continuity equations in the atmosphere. He was the first to find that these equations were "stiff" and were not solvable efficiently by any numerical technique then available. His efforts^{1, 2, 3} led to a computer simulation of the composition of the upper atmosphere that could be run inexpensively on the large scale computers available in those days. After these computer codes were applied to the undisturbed atmosphere⁴ and to several natural⁵ and manmade atmospheric disturbances, it became apparent that they could not predict long term atmospheric behavior.

(Received for publication 1 March 1984)

1. Keneshea, T.J. (1962) A Computer Program for Solving the Reaction Rate Equations in the E Ionospheric Region, AFCRL-62-828.
2. Keneshea, T.J. (1963) A Solution to the Reaction Rate Equations in the Atmosphere Below 150 Kilometers, AFCRL-63-711.
3. Keneshea, T.J. (1967) A Technique for Solving the General Reaction Rate Equations in the Atmosphere, AFCRL-67-0221, AD654010.
4. Keneshea, T.J., Narcisi, R.S., and Swider, W., Jr. (1970) Diurnal model of the E-region, J. Geophys. Res. 75:845-854.
5. Keneshea, T.J., and Fowler, R.J. (1966) Computed Electron, Ion, and Neutral Density Profiles for the Solar Eclipse of 12 November 1966, AFCRL-66-741, AD646975.

Such simple computer codes, in which only chemical reactions among the atmospheric gases are considered, are excellent for determining short term variations of the atmosphere, but give completely erroneous results for solutions beyond a few hours. In fact, these codes are valid only for solution times shorter than the transport time constants for the various atmospheric gases.

It then became obvious that any realistic atmospheric simulation should be based upon the equations of motion for the particles as well as their continuity equations. This report describes in detail a computer scheme for solving large sets of coupled partial differential equations of conservation of mass and momentum and also presents the computer code itself with instructions for using it. Also included are the results of an application of the code to the development of a one-dimensional midlatitude winter simulation of the composition of the mesosphere and lower thermosphere.

2. THE EQUATIONS OF STATE

The equation for the conservation of particles of a particular gas within a multicomponent gas mixture is

$$\frac{\partial n_i}{\partial t} = P_i - n_i L_i - \nabla \cdot \vec{\phi}_i \quad (1)$$

where n_i is the concentration of species i , P_i is the rate at which it is formed, and $n_i L_i$ is the rate at which it is removed through gas phase chemical reactions with other members of the ensemble. The vector flux of the species, $\vec{\phi}_i$, is induced by a combination of all nonchemical processes. In this report, only molecular and eddy diffusion will be considered in the flux term

$$\vec{\phi}_i = n_i \vec{C}_i + n_i \vec{V}_i \quad (2)$$

where \vec{C}_i is the velocity of species i resulting from molecular diffusion and \vec{V}_i is the velocity resulting from eddy diffusion. Since only the vertical distribution of species is to be considered here, the continuity equation can be written as

$$\frac{\partial n_i}{\partial t} = P_i - n_i L_i - \frac{\partial}{\partial Z} (n_i w_i + n_i v_i) \quad (3)$$

where w_i and v_i are the vertical components of the velocities \vec{C}_i and \vec{V}_i respectively.

The equation of motion for a neutral gas particle in a multicomponent gas mixture can be written as⁶

$$\frac{\partial \vec{C}_i}{\partial t} + \frac{1}{\rho_i} \nabla p_i + \frac{\alpha_i p_i}{\rho_i} \nabla \ln T - \vec{g} = - \sum_j \nu_{ij} (\vec{C}_i - \vec{C}_j) \quad (4)$$

where $\rho_i = n_i m_i$ is the mass density of species i , α_i is its thermal diffusion factor in a multicomponent mixture, p_i is its partial pressure, T is the temperature of the mixture, \vec{g} is the acceleration of gravity, and ν_{ij} is the collision frequency of species i with species j . The equation of motion is written here in terms of the collision frequencies between particles rather than the molecular diffusion coefficients in a multicomponent gas as was done by Keneshea and Zimmerman⁷ and Keneshea et al.⁸ The pressure of the gas is simply

$$p_i = n_i kT \quad (5)$$

where k is Boltzmann's constant. Therefore, in the one-dimensional case

$$\frac{1}{\rho_i} \nabla p_i + \alpha_i \frac{p_i}{\rho_i} \nabla \ln T + g = \frac{kT}{m_i} \left[\frac{1}{n_i} \frac{\partial n_i}{\partial Z} + \frac{(1 + \alpha_i)}{T} \frac{\partial T}{\partial Z} + \frac{1}{H_i} \right] \quad (6)$$

where m_i is the mass of species i and $H_i = kT/m_i g$ is its scale height. The equation of motion in the vertical direction for a neutral gas particle can then be written as

$$\frac{\partial w_i}{\partial t} + \frac{kT}{m_i} \left[\frac{1}{n_i} \frac{\partial n_i}{\partial Z} + \frac{(1 + \alpha_i)}{T} \frac{\partial T}{\partial Z} + \frac{1}{H_i} \right] = - \sum_j \nu_{ij} (w_i - w_j). \quad (7)$$

If inertial effects are ignored, then the equation of motion for ions and electrons in the presence of both magnetic and electric fields are⁹

-
- 6. Chapman, S., and Cowling, T.G. (1970) The Mathematical Theory of Non-Uniform Gases, Cambridge University Press
 - 7. Keneshea, T.J., and Zimmerman, S.P. (1970) The effect of mixing upon atomic and molecular oxygen in the 70-170 km region of the atmosphere, J. Atmos. Sci. 27:831-840.
 - 8. Keneshea, T.J., Zimmerman, S.P., and Philbrick, C.R. (1979) A dynamic model of the mesosphere and lower thermosphere, Planet. Space Sci. 27:385-401.
 - 9. Banks, P.M., and Kockarts, G. (1973) Aeronomy, Part A and Part B, Academic Press, New York.

$$\frac{\partial \vec{C}_i}{\partial t} + \frac{1}{\rho_i} \nabla p_i - \vec{g} - \frac{q_i}{m_i} (\vec{E} + \frac{1}{c} \vec{C}_i \times \vec{B}) = - \sum_j \nu_{ij} (\vec{C}_i - \vec{C}_j) \quad (8)$$

and

$$\frac{\partial \vec{C}_e}{\partial t} + \frac{1}{\rho_e} \nabla p_e - \vec{g} - \frac{q_e}{m_e} (\vec{E} + \frac{1}{c} \vec{C}_e \times \vec{B}) = - \sum_j \nu_{ej} (\vec{C}_e - \vec{C}_j)$$

where q is the particle charge, m is its mass, \vec{E} is the total electric field vector, c is the speed of light, and \vec{B} is the magnetic field vector. After much algebra, ignoring the horizontal component of the electric field as well as the horizontal velocities of the particles, and assuming that the electron and the ion temperatures are the same as the temperature of the neutral gas mixture, the equation of motion for the vertical velocities of the ionic species becomes

$$\begin{aligned} \frac{\partial w_i}{\partial t} + \frac{kT}{m_i} \left[\frac{1}{n_i} \frac{\partial n_i}{\partial Z} + \frac{1}{n_e} \frac{\partial n_e}{\partial Z} + \frac{2}{T} \frac{\partial T}{\partial Z} + \frac{1}{H_i} \right] \\ = -w_i \sum_j \nu_{ij} \left\{ \frac{\Omega_i^2 + \left(\sum_j \nu_{ij} \right)^2}{\Omega_i^2 \sin^2 I + \left(\sum_j \nu_{ij} \right)^2} \right\} + \sum_j \nu_{ij} w_j \end{aligned} \quad (9)$$

where n_e is the concentration of the electrons, I is the magnetic dip angle, and

$$\Omega_i = \frac{B_0 e}{m_i c} \quad (10)$$

where B_0 is the magnitude of the magnetic field and e is the charge on the electron.

The eddy diffusion velocity, taken from the formulation of Colegrove et al,¹⁰ is

$$v_i = -K \left[\frac{1}{n_i} \frac{\partial n_i}{\partial Z} + \frac{1}{T} \frac{\partial T}{\partial Z} + \frac{1}{H} \right] \quad (11)$$

where K is the vertical eddy diffusion coefficient and \bar{H} is the scale height of the mean mass.

10. Colegrove, F.D., Johnson, F.S., and Hanson, W.B. (1966) Atmospheric composition in the lower thermosphere, J. Geophys. Res. 71:2227-2236.

Whenever a photodissociation rate coefficient is needed in the calculations, the O₂ and O₃ column densities are estimated and a simple table look-up is made for the particular J_j required.

A short discussion of each photodissociation process and the photodissociation rate coefficient tables are presented in Appendix B.

6.2 The Photoionization Rate Coefficients

Reactions 185 through 191 and reactions 206 and 207 in the list of reactions in Appendix A are the photoionization processes considered in this report. The EUV solar fluxes, the total absorption cross sections for O₂, N₂, and O, and the ionization cross sections for O₂, N₂, O, H, and He are listed in Table B21. The EUV solar fluxes from 1 to 121.57 nm are from the compilation of Manson.¹⁸ The X-ray fluxes below 1 nm are taken from Swider.¹⁹ The cross sections for total absorption and for total photoionization of O₂, N₂, and O were obtained from a consideration of reported experimental and theoretical values. References, definitions, and a discussion of the absorption spectra can be found in Huffman²⁰ for wavelengths to 30.4 nm. For shorter wavelengths, the values given by Swider¹⁹ were used. A discussion of some photoionization rate coefficients can be found in Keneshea and Huffman.²¹ The photoionization cross sections for N, H, and He are taken from Banks and Kockarts.⁹

The photoionization of the metastable O₂(¹Δg) molecule has been included in these calculations. The wavelength region of interest is from 102.7 to 111.8 nm, where ground state O₂ is the primary absorber of the incident solar radiation but is not ionized itself. The effective absorption coefficient for O₂ and the ionization cross section for O₂(¹Δg) are taken from Huffman et al.²² Nitric oxide is ionized readily by Lyman α radiation. The ionization cross section for NO at Lyman α is 2 x 10⁻¹⁸ cm² as given by Watanabe.²³

-
18. Manson, J.E. (1976) Satellite Measurements of Solar UV During 1974, AFCRL-TR-76-0006, AD A021490.
 19. Swider, W., Jr. (1969) Ionization rates due to the attenuation of 1-100A non-flare solar X rays in the terrestrial atmosphere, Rev. Geophys. 7:573-594.
 20. Huffman, R.E. (1972) Photochemical processes. Cross-section data, Chapter 12, in Defense Nuclear Agency Reaction Rate Handbook, DNA 1948H.
 21. Keneshea, T.J., and Huffman, R.E. (1972) Solar Photoionization Rate Constants and Ultraviolet Intensities, AFCRL-72-0667.
 22. Huffman, R.E., Paulsen, D.E., Larrabee, J.C., and Cairns, R.B. (1971) Decrease in D-region O₂ (¹Δg) photoionization rates resulting from CO₂ absorption, J. Geophys. Res. 76:1028-1038.
 23. Watanabe, K. (1954) Photoionization and total absorption cross sections of gases. I. Ionization potentials of several molecules. Cross sections of NH₃ and NO, J. Chem. Phys. 22:1564-1570.

The total rate coefficient is obtained simply by summing $J_j(\lambda)$ over all the wavelengths.

$$J_j = \sum_{\lambda} \sigma'_j(\lambda) I_o(\lambda) e^{-\sum_i \sigma_i(\lambda) N_i}. \quad (54)$$

6.1 The Photodissociation Rate Coefficients

Reactions 192 through 205 in the list of reactions in Appendix A are the photodissociation processes considered in this report. In computing the transmittance, only molecular oxygen and ozone are considered as absorbers of the solar radiation. In addition to these two species, the absorption and subsequent photodissociation of H_2O , H_2O_2 , NO_2 , NO_3 , N_2O , HNO_2 , CO_2 , and HO_2 are discussed in Appendix B. Also presented in Appendix B are the photodissociation of N_2O_5 , HNO_3 , CH_4 , and CH_2O , even though these species are not included in the aeronomic calculations discussed in this report.

The incident solar flux $I_o(\lambda)$ and the absorption cross sections for O_2 , O_3 , H_2O , H_2O_2 , NO_2 , NO_3 , and N_2O used in calculating the photodissociation rate coefficients are listed in Table B1. The cross sections for O_2 in the wavelength region of the Schumann Runge bands (175-205 nm) are not tabulated here. In this wavelength region, the molecular oxygen dissociation data of Hudson and Mahle¹⁵ are used. The incident solar fluxes listed here are those tabulated by Ackerman.¹⁶ In the wavelength region 123 to 190 nm, however, the Ackerman fluxes have been replaced by the measurements of Heroux and Swirbalus.¹⁷ The absorption cross sections for N_2O_5 , HNO_2 , HNO_3 , CH_4 , CO_2 , CH_2O , and HO_2 are listed in Table B2.

In computing the time dependence of atmospheric composition, the computation of the photodissociation rate coefficients using Eq. (54) can be a time consuming process. As the solar zenith angle changes, the column densities N_i will change and, with them, the J_j 's. It is desirable to have available tables of these coefficients that are not dependent upon specific distributions of molecular oxygen and ozone. To keep the tables perfectly general, therefore, the coefficients are computed and tabulated in Appendix B as functions of the column densities of O_2 and O_3 .

15. Hudson, R.D., and Mahle, S.H. (1972) Photodissociation rates of molecular oxygen in the mesosphere and lower thermosphere, *J. Geophys. Res.* 77:2902-2914.

16. Ackerman, M. (1971) Ultraviolet solar radiation related to mesospheric processes, in *Mesospheric Models and Related Experiments*, G. Fiocco, Ed., D. Reidel, Dordrecht, Netherlands.

17. Heroux, L., and Swirbalus, R.A. (1976) Full-disc solar flares between 1230 and 1940 Å, *J. Geophys. Res.* 81:436-440.

the composition of the atmosphere is computed, the dissociation and ionization of atmospheric gases by incident solar radiation must be considered. In most calculations, these photoprocesses are handled as chemical reactions. Consequently, the rate coefficients for dissociation and ionization of the various gases must be known.

The photodissociation and photoionization rate coefficients for a gas particle at any altitude in the atmosphere is a product of the flux of the solar radiation arriving at that altitude and its photodissociation, or photoionization cross section. Several excellent reviews of the absorption properties of atmospheric gases can be found in the literature. Among these are Hudson¹² and Johnston and Graham.¹³ Also, an excellent survey article by Turco¹⁴ includes a discussion of photodissociation processes for several gases.

Beer's Law expresses the fractional transmission of radiation through a gas mixture

$$T(\lambda) = \frac{I(\lambda)}{I_0(\lambda)} = e^{-\tau(\lambda)} \quad (51)$$

where $I_0(\lambda)$ is the intensity of the incident radiation, $I(\lambda)$ is the intensity of the transmitted radiation, and $\tau(\lambda)$ is the optical depth of the absorbing layer, all at the wavelength λ . The optical depth is given by

$$\tau(\lambda) = \sum_i \sigma_i(\lambda) N_i \quad (52)$$

where $\sigma_i(\lambda)$ is the absorption cross section of the $i^{\text{'}}\text{th}$ type of particle and N_i is the total number of particles of this type in a cm^2 column along the ray path to the sun.

The rate coefficient for the photoprocess for j -type particles at wavelength λ is determined by

$$J_j(\lambda) = \sigma'_j(\lambda) I_0(\lambda) T(\lambda) \quad (53)$$

where $\sigma'_j(\lambda)$ is the cross section for photodissociation or photoionization of the gas.

- 12. Hudson, R.D. (1971) Critical review of ultraviolet photoabsorption cross sections for molecules of astrophysical and aeronomical interest, Rev. Geophys. Space Phys. 9:305-406.
- 13. Johnston, H.S., and Graham, R. (1974) Photochemistry of NO_x and HNO_x compounds, Can. J. Chem. 52:1415-1423.
- 14. Turco, R.P. (1975) Photodissociation rates in the atmosphere below 100 km, Geophys. Surveys 2:153-192.

Therefore, the average collision frequency for charge exchange is

$$\bar{\nu}_{in} = 1.73725 n \gamma. \quad (48)$$

The average reduced charge exchange collision frequencies between like particles are given in Table 2. Again, it is assumed that the ion and neutral gas temperatures are equal.

Table 2. The Average Reduced Charge Exchange Collision Frequencies Between Like Particles

$H^+ + H$	$1.0 \times 10^{-10} H (2T)^{1/2}$
$O^+ + O$	$1.6 \times 10^{-11} O (2T)^{1/2}$
$N^+ + N$	$1.7 \times 10^{-11} N (2T)^{1/2}$
$He^+ + He$	$3.0 \times 10^{-11} He (2T)^{1/2}$
$O_2^+ + O_2$	$1.1 \times 10^{-11} O_2 (2T)^{1/2}$
$N_2^+ + N_2$	$2.2 \times 10^{-11} N_2 (2T)^{1/2}$

The average ion-ion momentum transfer collision frequency is given by

$$\bar{\nu}_{ij} = 1.3 n_j \left(\frac{1}{\mu^{1/2} T^{3/2}} \right) \text{sec}^{-1} \quad (49)$$

assuming that the charge states of the colliding particles are both unity.

The collision frequency between i and j-type particles is then simply

$$\nu_{ij} = \frac{m_i}{(m_i + m_j)} \bar{\nu}_{ij}. \quad (50)$$

6. THE CHEMISTRY AND PHOTOCHEMISTRY

The chemical reactions included in this model are listed in Appendix A. When

Table 1. Polarizabilities of Neutral Gases

GAS	$\alpha(10^{-24} \text{ cm}^3)$
N ₂	1.76
O ₂	1.59
O ₂ (¹ Δg)	1.59
O	0.79
H	0.667
H ₂	0.82
N	1.1
N(² D)	1.1
NO	1.74
He	0.21
N ₂ O	3.00
CO ₂	2.63
CO	1.97

Let

$$\bar{Q}_E = \sigma = \frac{\gamma}{v} \quad (46)$$

where γ is the chemical reaction rate constant and v is the velocity of the colliding particle.

Since

$$v = \left(\frac{3kT}{m}\right)^{1/2}, \text{ then}$$

$$\bar{v}_{in} = \frac{4}{3} n \left(\frac{16kT}{\pi m}\right)^{1/2} \frac{\gamma}{\left(\frac{3kT}{m}\right)^{1/2}}. \quad (47)$$

5. COLLISION FREQUENCIES

A good review of collision processes can be found in Banks and Kockarts.⁹
All of the material that follows was taken from that source.

The average neutral-neutral collision frequencies are given by

$$\bar{\nu}_{ij} = 2n_j \sigma_{ij}^2 \left(\frac{2\pi kT}{\mu} \right)^{1/2} \text{ sec}^{-1} \quad (43)$$

where μ , the reduced mass, is given by

$$\mu = \frac{m_i m_j}{m_i + m_j}$$

and

$$\sigma_{ij} = \frac{1}{2}(\sigma_i + \sigma_j)$$

where σ_i and σ_j are the diameters of the colliding particles.

The average collision frequency for momentum transfer between ions and neutral particles is

$$\bar{\nu}_{in} = 2.6 \times 10^{-9} n_n \left(\frac{\alpha}{\mu} \right)^{1/2} \text{ sec}^{-1} \quad (44)$$

where n_n is the concentration of the neutral particles, α is the neutral gas atomic polarizability in units of 10^{-24} cm^3 , and μ is in amu. Table 1 lists the polarizability of several atmospheric gases.⁹

The average collision frequency for charge exchange between unlike particles is given by

$$\bar{\nu}_{in} = \frac{4}{3} n \left(\frac{8k}{\pi m} \right)^{1/2} (2T)^{1/2} \bar{Q}_E \quad (45)$$

where \bar{Q}_E is the average charge transfer cross section for ion-neutral collisions. In Eq. (45), the ion and neutral temperatures are assumed to be equal.

The calculation of F_K , however, requires the upper boundary velocity at the end of the time step. At the upper boundary, O_2^+ , NO^+ , O_3 , OH , H , HO_2 , H_2O , H_2O_2 , H_2 , NO_2 , O_2 (Δg), and argon are all minor species. It is assumed, therefore, that they are in diffusive equilibrium there. This means that their diffusion velocities are zero at the upper boundary, but they are still subject to the mean mass motion.

The remaining species, however, do have diffusion velocities at the upper boundary that must be defined. It is impossible to calculate the upper boundary velocities from Eq. (18) because concentrations are required at $K + 1$. Since this altitude is outside the volume, the concentrations are not known there. Therefore, to estimate the velocities at the upper boundary at the end of the time step, we investigate the continuity equation, Eq. (3), at the upper boundary where $v_i = 0$ (since the upper boundary is outside of the turbulence region).

$$\frac{\partial n_i}{\partial t} = P_i - n_i L_i - \frac{\partial}{\partial Z} (n_i w_i). \quad (41)$$

Setting the time derivative to zero and solving the resulting equation for dw_i/dZ gives

$$\frac{dw_i}{dZ} = - \frac{w_i}{n_i} \frac{dn_i}{dZ} + \frac{P_i - n_i L_i}{n_i}. \quad (42)$$

This ordinary differential equation can be solved numerically over the last height step to give values of w_i at the upper boundary. Since Eq. (42) is solved in every iteration, the final iteration will give the upper boundary value of w_i at the end of the time step.

The new concentration profiles at $t + \Delta t$ are therefore computed as follows: Starting at the lower boundary, the E and F arrays are computed in the increasing height direction up to the upper boundary. The new concentrations are then computed starting at the upper boundary and applying Eq. (27) in the decreasing altitude direction down to the lower boundary. This procedure is repeated over each time step until the solutions converge for that time step. Once a solution has been accepted, the time step is doubled (provided it has not reached the maximum value set by the program). Using the concentrations and velocities computed at the end of the last time step, the equations are solved over the next time step. In this manner, the solution advances.

Eq. (36) can be written as

$$\tilde{B}_K^\ell n_K^{\ell+1} - \tilde{C}_K^\ell n_{K-1}^{\ell+1} = \tilde{D}_K^\ell \quad (37)$$

where

$$\tilde{A}_K^\ell = 0$$

$$\tilde{B}_K^\ell = -A_K^\ell \left[1 - \Gamma_K \Delta Z_K - \frac{m_i \Delta Z_K}{k T_K} \sum_j \nu_{ij} (w_i - w_j)_K^{\ell+1} \right] + B_K^\ell$$

$$\tilde{C}_K^\ell = C_K^\ell$$

$$\tilde{D}_K^\ell = D_K^\ell. \quad (38)$$

Eq. (28) can be rewritten at level K as

$$E_K = \frac{\tilde{A}_K^\ell}{[\tilde{B}_K^\ell - \tilde{C}_K^\ell E_{K-1}]}$$

and

$$F_K = \frac{\tilde{D}_K^\ell + \tilde{C}_K^\ell F_{K-1}}{\tilde{B}_K^\ell - \tilde{C}_K^\ell E_{K-1}}. \quad (39)$$

It is obvious, then, that

$$E_K = 0$$

and

$$F_K = \frac{D_K^\ell + C_K^\ell F_{K-1}}{B_K^\ell - A_K^\ell \left[1 - \Gamma_K \Delta Z_K - \frac{m_i \Delta Z_K}{k T_K} \sum_j \nu_{ij} (w_i - w_j)_K^{\ell+1} \right] - C_K^\ell E_{K-1}}. \quad (40)$$

$$n_o^{\ell+1} = \frac{n_o^\ell + P_o^{\ell+1} \Delta t}{[1 + L_o^{\ell+1} \Delta t]}.$$
(30)

E_o and F_o at the lower boundary are simply

$$E_o = 0$$

and

$$F_o = n_o^{\ell+1}.$$
(31)

At the upper boundary, the time derivative of the velocities in Eq. (7) is set to zero:

$$\frac{kT}{m_i} \left[\frac{1}{n_i} \frac{dn_i}{dZ} + \frac{(1+\alpha_i)}{T} \frac{dT}{dZ} + \frac{1}{H_i} \right] = - \sum_j \nu_{ij} (w_i - w_j).$$
(32)

Solving Eq. (32) for dn_i/dZ gives

$$\frac{dn_i}{dZ} = -n_i \Gamma_i - \frac{n_i m_i}{kT} \sum_j \nu_{ij} (w_i - w_j).$$
(33)

Writing Eq. (33) at the upper boundary (K) in finite difference notation gives

$$\frac{n_{K+1}^{\ell+1} - n_K^{\ell+1}}{\Delta Z_K} = n_K^{\ell+1} \Gamma_K - \frac{n_K^{\ell+1} m_i}{kT_K} \sum_j \nu_{ij} (w_i - w_j)_K^{\ell+1}.$$
(34)

Solving Eq. (34) for $n_{K+1}^{\ell+1}$ gives

$$n_{K+1}^{\ell+1} = n_K^{\ell+1} \left[1 - \Gamma_K \Delta Z_K - \frac{m_i \Delta Z_K}{kT_K} \sum_j \nu_{ij} (w_i - w_j)_K^{\ell+1} \right].$$
(35)

Replacing $n_{K+1}^{\ell+1}$ in Eq. (25) written at level K with Eq. (35) gives

$$\left\{ -A_K^\ell \left[1 - \Gamma_K \Delta Z_K - \frac{m_i \Delta Z_K}{kT_K} \sum_j \nu_{ij} (w_i - w_j)_K^{\ell+1} \right] + B_K^\ell \right\}$$

$$n_K^{\ell+1} - C_K^\ell n_{K-1}^{\ell+1} = D_K^\ell$$
(36)

From Eqs. (26) and (27), the following definitions are obtained:

$$E_k = \frac{A_k^{\ell}}{[B_k^{\ell} - C_k^{\ell} E_{k-1}]} \quad (27)$$

and

$$F_k = \frac{D_k^{\ell} + C_k^{\ell} F_{k-1}}{[B_k^{\ell} - C_k^{\ell} E_{k-1}]} \quad (28)$$

Using Eqs. (27) and (28), the height profiles for the concentrations can be calculated provided $n_O^{\ell+1}$ and $n_K^{\ell+1}$ can be specified at the lower and upper boundaries respectively. The boundary conditions will be discussed in Section 4.

It should be noted that only some of the terms on the right-hand side of the equations for A_k^{ℓ} , B_k^{ℓ} , C_k^{ℓ} , and D_k^{ℓ} are specified at the beginning of the time step. The chemistry terms and the velocity sums are specified at the end of the time step. Because of this, the solutions require iteration over the time step. At the end of each iteration of the concentrations, the chemistry and the velocities from Eq. (18) are recomputed. New concentrations are then calculated using these updated values. This procedure is repeated until, on successive iterations, the concentrations of all species change by less than 1 percent at all altitudes.

4. BOUNDARY CONDITIONS

At the lower boundary, the concentrations of N_2 , O_2 , H_2O , H_2 , and Ar are held fixed for all time at their initial values ($n_O^{\ell+1} = \text{constant}$). All of the other species in the transport mode prove to be chemically controlled at 50 km, so their concentrations are computed from the simplified continuity equations. With the divergence of the flux set to zero in Eq. (1), that equation becomes

$$\frac{dn_i}{dt} = P_i - n_i L_i \quad (29)$$

If the derivative in Eq. (29) is replaced by simple differences, then, at the lower boundary, we have

in which

$$Q_k = \frac{1}{[1 + \Delta t \lambda \sum_j \nu_{ij}]}$$

$$S_k = (1 - \Gamma_k \Delta Z_k)$$

$$\bar{S}_k = (1 - \bar{\Gamma}_k \Delta Z_k).$$

It is possible, therefore, to write a tridiagonal matrix between the lower and upper boundaries for each species using Eq. (21). These tridiagonal matrices are then solved for the vertical distribution of the concentration using the method of Richtmyer and Morton.¹¹

Substituting the recurrence relation

$$n_k^{\ell+1} = E_{k-1} n_k^{\ell+1} + F_{k-1} \quad (23)$$

into Eq. (21) gives

$$-A_k^\ell n_{k+1}^{\ell+1} + B_k^\ell n_k^{\ell+1} - C_k^\ell E_{k-1} n_k^{\ell+1} - C_k^\ell F_{k-1} = D_k^\ell. \quad (24)$$

This can be written as

$$-A_k^\ell n_{k+1}^{\ell+1} + [B_k^\ell - C_k^\ell E_{k-1}] n_k^{\ell+1} = D_k^\ell + C_k^\ell F_{k-1}. \quad (25)$$

Solving Eq. (25) for $n_k^{\ell+1}$ gives

$$n_k^{\ell+1} = \frac{A_k^\ell}{[B_k^\ell - C_k^\ell E_{k-1}]} n_{k+1}^{\ell+1} + \frac{D_k^\ell + C_k^\ell F_{k-1}}{[B_k^\ell - C_k^\ell E_{k-1}]} . \quad (26)$$

Eq. (23) written at level k is

$$n_k^{\ell+1} = E_k n_{k+1}^{\ell+1} + F_k. \quad (27)$$

11. Richtmyer, R.D., and Morton, K.W. (1967) Difference Methods for Initial-Value Problems, Interscience Publishers, New York.

$$\begin{aligned}
\phi_{k-1}^{\ell+1} = & \frac{1}{\left[1 + \Delta t \lambda_i \sum_j \nu_{ij}\right]} \left\{ n_{k-1}^{\ell+1} w_{k-1}^{\ell+1} - \frac{k T_{k-1}}{m_i} \frac{\Delta t}{\Delta Z_{k-1}} \right. \\
& \left. \left[n_k^{\ell+1} - n_{k-1}^{\ell+1} (1 - \bar{F}_{k-1} \Delta Z_{k-1}) \right] \right. \\
& \left. + n_{k-1}^{\ell+1} \Delta t \sum_{j \neq i} \nu_{ij} w_{j_{k-1}}^{\ell+1} \right\} \\
& - \frac{K_{k-1}}{\Delta Z_{k-1}} \left\{ n_k^{\ell+1} - n_{k-1}^{\ell+1} (1 - \bar{F}_{k-1} \Delta Z_{k-1}) \right\}. \tag{20}
\end{aligned}$$

Substituting Eqs. (19) and (20) into Eq. (12) and collecting like terms in n gives

$$- A_k^\ell n_{k+1}^{\ell+1} + B_k^\ell n_k^{\ell+1} - C_k^\ell n_{k-1}^{\ell+1} = D_k^\ell \tag{21}$$

where

$$\begin{aligned}
A_k^\ell &= \frac{\Delta t}{\Delta Z_k \Delta Z_{k-1}} \left[Q_k \frac{k T_k \Delta t}{m} + K_k \right] \\
B_k^\ell &= 1 + L_k^{\ell+1} \Delta t + \frac{\Delta t}{\Delta Z_k \Delta Z_{k-1}} \\
&\left[Q_k \left(w_k^\ell \Delta Z_k + S_k \frac{k T_k \Delta t}{m} + \Delta t \Delta Z_k \sum_{j \neq i} \nu_{ij} w_k^{\ell+1} \right) \right. \\
&\left. + \bar{S}_k K_k + Q_{k-1} \frac{k T_{k-1} \Delta t}{m} \frac{\Delta Z_k}{\Delta Z_{k-1}} + K_{k-1} \frac{\Delta Z_k}{\Delta Z_{k-1}} \right] \\
C_k^\ell &= \frac{\Delta t}{\Delta Z_k \Delta Z_{k-1}} \left[Q_{k-1} \left(w_{k-1}^\ell \Delta Z_k + S_{k-1} \frac{k T_{k-1} \Delta t}{m} \frac{\Delta Z_k}{\Delta Z_{k-1}} \right. \right. \\
&\left. \left. + \Delta t \Delta Z_k \sum_{j \neq i} \nu_{ij} w_{k-1}^{\ell+1} \right) + \bar{S}_k K_{k-1} \frac{\Delta Z_k}{\Delta Z_{k-1}} \right] \\
D_k^\ell &= n_k^\ell + P_k^{\ell+1} \Delta t \tag{22}
\end{aligned}$$

both computed at level k. Also

$$\lambda_i = \frac{\Omega_i^2 + \left(\sum_j v_{ij} \right)^2}{\Omega_i^2 \sin^2 I + \left(\sum_j v_{ij} \right)^2} \quad (17)$$

The finite difference equations, Eqs. (13) and (14), can be solved for the velocities $w_k^{\ell+1}$ at the end of the time step:

$$w_k^{\ell+1} = \frac{1}{\left[1 + \Delta t \lambda_i \sum_j v_{ij} \right]} \left\{ w_k^\ell - \frac{kT_k}{m_i} \frac{\Delta t}{\Delta Z_k} \left[\frac{n_{k+1}^{\ell+1} - n_k^{\ell+1}}{n_k^{\ell+1}} + \Gamma_k \Delta Z_k \right] + \Delta t \sum_j v_{ij} w_{jk}^{\ell+1} \right\} \quad (18)$$

This equation holds for both the ionic and neutral species using the above definitions of Γ_k and with $\lambda_i = 1$ for the neutral species. The total molecular and eddy diffusion flux at level k can then be written as

$$n_k^{\ell+1} w_k^{\ell+1} = \Phi_k^{\ell+1} = \frac{1}{\left[1 + \Delta t \lambda_i \sum_j v_{ij} \right]} \left\{ n_{k+1}^{\ell+1} w_k^\ell - \frac{kT_k}{m_i} \frac{\Delta t}{\Delta Z_k} \left[n_{k+1}^{\ell+1} - n_k^{\ell+1} \left(1 - \Gamma_k \Delta Z_k \right) \right] + n_k^{\ell+1} \Delta t \sum_{j \neq i} v_{ij} w_{jk}^{\ell+1} \right\} - \frac{K_k}{\Delta Z_k} \left\{ n_{k+1}^{\ell+1} - n_k^{\ell+1} \left(1 - \bar{\Gamma}_k \Delta Z_k \right) \right\} \quad (19)$$

where

$$\bar{\Gamma}_k = \frac{1}{T_k} \frac{\partial T}{\partial Z} + \frac{1}{H} \quad \text{and} \quad \bar{H} = \frac{kT}{mg}.$$

A similar expression can be derived for $\Phi_{k-1}^{\ell+1}$ at level k-1.

3. THE FINITE DIFFERENCE EQUATIONS

If l denotes the grids on the time mesh and k the grids on the height mesh, then the continuity equations, Eq. (1), can be written in finite difference notation at the height level k as

$$\frac{n_k^{l+1} - n_k^l}{\Delta t} = P_k^{l+1} - n_k^{l+1} L_k^{l+1} - \frac{\phi_k^{l+1} - \phi_{k-1}^{l+1}}{\Delta Z_{k-1}} \quad (12)$$

where, for the sake of clarity, the subscript i has been dropped. Using the equations of motion, Eqs. (7) and (9), the ϕ_k^{l+1} and ϕ_{k-1}^{l+1} will be replaced in Eq. (12) with their equivalent expressions.

The equations of motion, Eq. (7), for the neutral species can be written in finite difference form at height level k as

$$\begin{aligned} \frac{w_k^{l+1} - w_k^l}{\Delta t} = & - \frac{k T_k}{m_i} \left[\frac{n_{k+1}^{l+1} - n_k^{l+1}}{n_k^{l+1} \Delta Z_k} + \Gamma_k \right] - w_k^{l+1} \sum_j \nu_{ij} \\ & + \sum_{j \neq i} \nu_{ij} w_{jk}^{l+1} \end{aligned} \quad (13)$$

and Eq. (9) for the ionic species as

$$\begin{aligned} \frac{w_k^{l+1} - w_k^l}{\Delta t} = & - \frac{k T_k}{m_i} \left[\frac{n_{k+1}^{l+1} - n_k^{l+1}}{n_k^{l+1} \Delta Z_k} + \Gamma_k \right] - w_k^{l+1} \lambda_i \sum_j \nu_{ij} \\ & + \sum_{j \neq i} \nu_{ij} w_{jk}^{l+1}. \end{aligned} \quad (14)$$

For the neutral species,

$$\Gamma_k = \frac{(1 + \alpha_i)}{T} \frac{\partial T}{\partial Z} + \frac{1}{H_i} \quad (15)$$

and for the ionic species

$$\Gamma_k = \frac{1}{n_e} \frac{\partial n_e}{\partial Z} + \frac{2}{T} \frac{\partial T}{\partial Z} + \frac{1}{H_i} \quad (16)$$

In computing the transmittance of EUV radiation, O, O₂, and N₂ are considered as the primary absorbers. The photoionization rate coefficients are computed by using Eq. (54), where the sum is taken from 0.1 to 102.57 nm. Radiation below 15 nm is considered to be X-rays, and its ionization is partitioned among O, O₂, N₂, and N in accordance with Swider.¹⁹

The calculations also take into account the scattering of some radiation from the sunlit hemisphere into the nighttime hemisphere. These scattered radiations include Lyman α at 121.57 nm, Lyman β at 102.57 nm, the HE I line at 58.4 nm, and the HE II line at 30.4 nm. For simplicity, the amount of radiation that scatters into the dark hemisphere is taken to be a fraction of the direct noontime radiation at the particular altitude under consideration. These fractions are:

Lyman α , 0.01; Lyman β , 0.004; HE I, 0.0011; and HE II, 0.001.

Finally, the contribution of cosmic rays to the photoionization rate coefficients is also included. This is simply²⁴

$$q_i = 1 \times 10^{-17} n_i. \quad (55)$$

6.3 Electron Impact Ionization

Reactions 208 through 215 in the list of reactions in Appendix A are the energetic electron processes. The flux of energetic electrons appropriate to the calculations made here were supplied by Jasperse.²⁵ The ionization rate coefficients for the energetic electron reactions are computed from⁹

$$I = \int_{\xi} \sigma(\xi) \Phi(\xi) d\xi \quad (56)$$

where $\sigma(\xi)$ is the electron impact ionization cross section and $\Phi(\xi)$ is the energetic electron flux, both at energy ξ . The ionization cross section is given by⁹

$$\sigma(\xi) = \int_{\frac{\xi+I}{2}}^{\frac{\xi+I}{2}} S(\xi, W) dW \quad (57)$$

24. Nicolet, M., and Aikin, A.C. (1960) The formation of the D region of the ionosphere, *J. Geophys. Res.* 65:1469-1483.

25. Jasperse, J.R. (1978) Private communication.

in which

$$S(\xi, W) = \left(\frac{q_0 c_0 f_0}{W^2} \right) \left(\frac{I}{W} \right)^P \left[1 - \left(\frac{W}{E} \right)^\gamma \right]^\nu \left(\frac{W}{\xi} \right)^\Omega$$

where I is the ionization threshold for a particular ion state, $q_0 = 6.51 \times 10^{-14}$ ev² cm², $c_0 f_0$ is a constant, W is the excitation threshold energy in ev, and P , γ , ν , and Ω are parameters obtained from the best fit of Eq. (58) to experimental and theoretical results. Table B22 lists the values of these parameters used in the model presented here. The values of the rate coefficients computed for the energetic electron reactions are given in Table B23.

6.4 The Thermal Diffusion Factors

The thermal diffusion factors are assumed to be zero except for the following species: O(-0.27), O₂(+0.08), H(-0.39), H₂(-0.31), and He(-0.36). A discussion of the thermal diffusion factors can be found in Zimmerman and Keneshea,²⁶ from which source these values were taken.

7. A MID-LATITUDE ATMOSPHERE SIMULATION

The computer code presented here describes the distribution of 56 individual atmospheric species. Not all of them have the full coupled sets of equations in the system, however. Some gases in the atmosphere have such short chemical time constants that they will react chemically long before they can be moved out of a unit volume by any transport process. The vertical distribution of these species can be determined quite accurately by considering only the chemical reactions in which they are involved and ignoring any transport effects on them.

The species for which the coupled sets of mass and momentum equations are solved are:

O	H	N	CO	AR	H ⁺
O ₂	H ₂	N(² D)	CO ₂		HE ⁺
O ₂ (¹ A _g)	HO ₂	N ₂			O ⁺
O ₃	H ₂ O	NO			O ₂ ⁺
OH	H ₂ O ₂	NO ₂			NO ⁺
		N ₂ O			N ⁺

Those species whose distributions are determined neglecting transport are

²⁶. Zimmerman, S. P., and Keneshea, T.J. (1976) The thermosphere in motion, *J. Geophys. Res.* 81:3187-3197.

$O(^1D)$, N_2^+ , $O(^2D)$, and the electrons. The height distributions of all 56 species are computed over the entire altitude range, which is from 50 to 400 km.

In addition, the vertical distributions of the following species are computed over a reduced altitude range. The temporal variation of these species is also computed with the transport term removed from their continuity equations. The distributions of these species are computed between 50 and 120 km because, above this altitude, they do not exist in significant quantities. These species are:

O^-	O_4^+	$H_5O_2^+ \cdot CO_2$	HNO_2
O_2^-	NO_2^+	$O_2^+ \cdot H_2O$	NO_3
O_3^-	$H^+ \cdot H_2O$	$NO^+ \cdot H_2O$	
O_4^-	$H^+ \cdot (H_2O)_2$	$NO^+ \cdot (H_2O)_2$	
NO_2^-	$H^+ \cdot (H_2O)_3$	$NO^+ \cdot (H_2O)_3$	
NO_3^-	$H^+ \cdot (H_2O)_4$	$NO^+ \cdot CO_2$	
CO_3^-	$H_3O^+ \cdot N_2$		
CO_4^-	$H_3O^+ \cdot OH$		
		$H_3O^+ \cdot CO_2$	

Neither negative cluster ions nor metallic ions are considered here. To ensure that the atmosphere remains electrically neutral, one of the negatively charged species is always computed from the imposed condition of balance of charge. Below 120 km, where there are several types of negative ions in addition to the electrons, the most abundant negatively charged species is always computed from conservation of charge. That is, its concentration is computed as the difference between the sum of all the positive and the sum of all the other negative ions. Above 120 km, where the only negatively charged particles are the electrons, their concentration is always set equal to the sum of all the positive ions.

7.1 Initial Conditions

Throughout the development of the time-dependent solution of Eq. (1), the temperatures and the turbulent diffusion coefficients are held fixed at all altitudes in the code presented here. For the simulation developed in this report, both of these parameters were derived from data obtained during the ALADDIN I experiments of

20 November 1970.^{27,28} The temperatures shown in Figure 1 between 50 and 125 km were taken from measurements of Theon and Horvath.²⁹ From 125 to 160 km, the measurements of Golomb²⁷ were used. Above 160 km, the temperatures were taken from the 900 K exospheric temperature winter model of the U.S. Standard Atmosphere Supplements, 1966. A discussion of the turbulent diffusion coefficients obtained from the ALADDIN I experiments can be found in Keneshea et al.⁸ The coefficients presented there represent the minimum turbulent diffusion coefficients determined from the observations on the chemical trails. Following arguments given there, the turbulent diffusion coefficients (Figure 2) used in these calculations are three times larger than those presented in that paper.

Keneshea et al.⁸ found that stability considerations required that the height steps in the mesosphere not exceed 150 m. In the thermosphere, however, stability could be maintained with somewhat larger height steps. Therefore, to reduce the amount of core storage required for saving the numerous arrays as well as to reduce the amount of computer time per solution, a variable height step array was used. From the lower boundary, 50 km, up to 90 km, the height steps were fixed at 150 m. From 90 km up to the upper boundary, 400 km, the height steps were computed by assuming that they were proportional to the pressure scale height of the atmosphere. This produced an altitude array with 692 grid points, with the maximum height step at the upper boundary of about 1.5 km.

The initial concentration and average velocity profiles were taken from the final solution of an unpublished simulation using turbulent diffusion coefficients one third smaller than those shown in Figure 2. A discussion of a method for establishing initial concentrations on all the species is given in Keneshea et al.⁸ for a situation where no previous information is available on the vertical distribution of the various species.

The solutions to Eqs. (1) and (18) are started at noon of solution day 1 with the above initial values. Because this is a real time-dependent calculation which takes into account the daily variation of the solar zenith angle, no truly steady state condition on the species concentrations is ever achieved. The calculations are continued, therefore, until all species reproduce their concentrations diurnally at all

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- 27. Rosenberg, N.W., Golomb, D., Zimmerman, S.P., Vickery, W.K., and Theon, J.S. (1973) The ALADDIN Experiment-Part 1, Dynamics, Space Research XIII, Akademie-Verlag, Berlin, 435-439.
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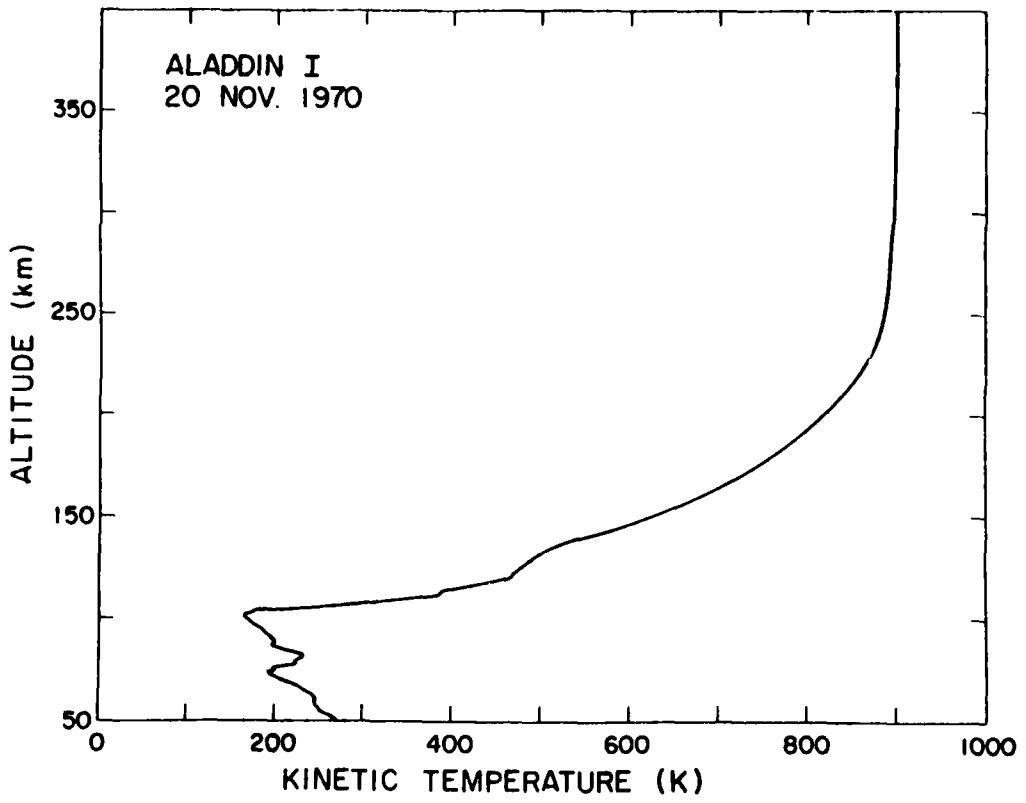


Figure 1. The Static Temperature Profile

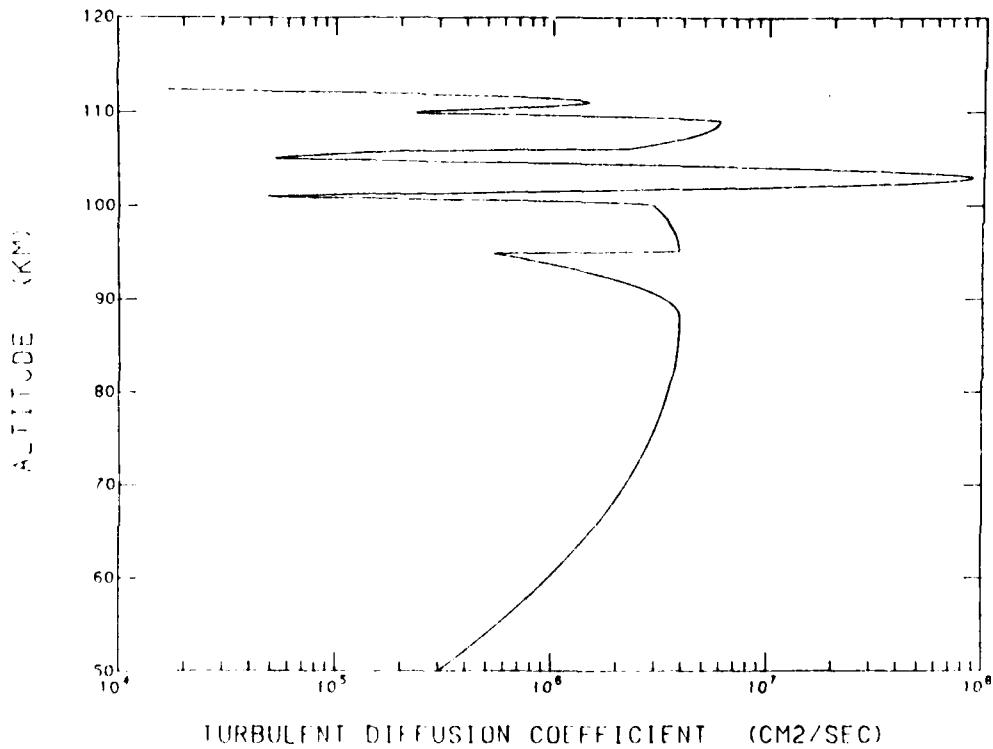


Figure 2. The Turbulent Diffusion Coefficients as Determined From the Chemical Release Data Obtained From the ALADDIN I Experiments

altitudes to within 1 percent. This condition of diurnal reproducibility is generally arrived at after about 20 solution days.

7.2 Thermospheric Results

The results to be discussed here were taken from the final diurnal cycle of the calculations when the major and most of the minor species at all altitudes were varying by no more than 1 percent from values achieved 24 hours earlier. The vertical distribution of the major species N_2 , O_2 , O, and argon are shown in Figure 3. In this one-dimensional temperature-independent calculation, N_2 , O_2 , and argon show no diurnal change, having arrived at an equilibrium with the invariant temperature given in Figure 1. Also displayed in Figure 3 is the fixed profile of helium used in these calculations.

Because of the constant temperature profile used here, atomic oxygen also shows no diurnal variation in the thermosphere down to about 120 km. Below this altitude, the chemistry dominates the production (dissociation of O_2) and loss of this species.⁷ Thus, we observe in Figures 3 and 4 large diurnal variations of atomic oxygen in the mesosphere.

Presented in Figure 5 are the mesosphere and lower thermosphere noontime profiles of the remaining important neutral species. Carbon dioxide is invariant in the mesosphere and lower thermosphere up to about 140 km. Above this altitude, the diurnal variation of the photodissociation process becomes important in the determination of the carbon dioxide concentrations. The variation of the CO_2/Ar ratio calculated here is shown in Figure 6 and compares well to that determined from atmospheric measurements of Trinks and Fricke.³⁰

Isodensity contour plots of the minor neutral species that exhibit a large diurnal variation are shown in Figures 7 to 11. The numbers in all the isodensity plots presented here are the common logarithms of the concentrations. These species, for the most part, are important in the determination of the electron density throughout the atmosphere. Above about 100 km, the charge distribution is dominated by NO^+ , O_2^+ , O^+ , He^+ , and H^+ . Thus, the resultant thermospheric charge balance follows the chain created primarily by the time-dependent distributions of the neutral species N, NO, and $N(^2D)$ as shown in Figures 7, 8, and 9, respectively. The other neutral thermospheric species in the code have little effect in determining the charge distribution.

The ion and electron isodensity contours determined from these neutrals are shown in Figures 12 to 19. As expected, the dominant positive ion in the upper

30. Trinks, H., and Fricke, K.H. (1978) Carbon dioxide concentrations in the lower thermosphere, J. Geophys. Res. 83:3883-3886.

thermosphere is the O^+ ion. The electron density distribution, therefore, follows the distribution of this ion very closely. Also displayed in Figure 19 is the altitude profile of $f_o F_2$, the space-time dependence of the peak of the electron density along with the measurements of this parameter by Evans.³¹ The correspondence is quite good, clearly demonstrating the midday altitude and amplitude and the night-time transition to higher altitudes and lower amplitudes. No horizontal winds are included in this one-dimensional simulation; yet, we obtain an excellent representation of the F_2 layer maximum.

Our rationale, then, is that the contribution of chemical production and loss coupled to the convergences and divergences created in our calculations are the main mechanisms that control the amplitude and altitude variation of the F_2 layer maximum. The question that arises from this one-dimensional calculation, which apparently gives good agreement with some F_2 layer maximum measurements, is: What is the effect of horizontal winds? The answer to this question may be inferred from measurements of Reinisch,³² who determined the time dependence of $f_o F_{max}$ for another period from ionosonde measurements. His results for $f_o F_{max}$ are shown in Figure 20. We observe a smaller peak ionization amplitude than Reinisch does, but, generally, the time dependence of his F_2 layer maximum follows our calculations quite well. Also observed are the large variations that presumably arise from local winds that show multireversals during the measuring period. The point we wish to stress is that ionization calculated from internally determined neutral species distributions does result in an internally consistent F_2 layer ionization behavior that compares well with some of the observed ionic behavior. An example of this is the recent work of Sethia et al.,³³ who have performed calculations similar to ours but include the effects of horizontal winds. They determine the distributions along magnetic field lines, while we calculate in the vertical direction with the magnetic field entering by way of the Larmor frequency and the magnetic dip angle. They input the neutral atmosphere, while we calculate it as an internally consistent integral part of the model.

In Figure 21, we have superimposed upon our calculation of the total electron content that taken from Figure 3, model 4 (no winds) of Sethia et al.³³ We observe that the shapes of the distributions are almost identical between the two one-dimensional models. The differences in magnitude can be attributed to the

31. Evans, J.V. (1971) Observations of F-region vertical velocities at Millstone Hill-I. Evidence for drifts due to expansion, contraction, and winds, Radio Sci. 6:609-626.

32. Reinisch, B. (1983) Private communication.

33. Sethia, G.C., Bailey, G.J., Moffett, R.J., and Hargreaves, J.K. (1983) The effects of neutral air winds on the electron content of the mid-latitude ionosphere and protonosphere in summer, Planet. Space Sci. 31:377-387.

different conditions under which the models were computed. We also compare (Figure 22) the total electron content to some measurements taken at Boulder by Sethia et al³³ as shown in their Figure 8 (July 2, no wind, and July 5, with wind). Again, the shape of our distribution compares well with the measured profiles. For those cases where the wind amplitude is not very large, the contribution of the $\vec{v} \times \vec{B}$ mechanism is very much smaller than that of the chemistry. We make another comparison with the work of Poulter et al³⁴ shown in Figure 23. Here we examine the total electron content for measurements at Boulder and Lancaster (Figure 5 of Poulter et al), and again quite good agreement is seen between the shape of the distributions from the chemical-transport one-dimensional model and the measurements. The observed differences between the Boulder and the Lancaster curves result from the differences in the raypath geometries. Comparison is also quite good between our calculations and the $F_o F_2$ measured by Kohl et al³⁵ and discussed by Risbeth.³⁶

We reiterate the point we wish to stress (see also Rush et al³⁷): A comprehensive, one-dimensional, time-dependent ion-neutral model demonstrates excellent agreement with measurements of electron densities. The inclusion of the effects of winds is necessary, however, to explain the deviation of the ions and/or electrons from their equilibrium distributions, which significantly vary over the diurnal cycle.

7.3 Mesospheric Results

The ionization in the mesosphere is much more complex than the ionization in the thermosphere. Ablating meteorites create long-lived atomic ions from the metallic species present in the atmosphere. These metallic ions are significant and are observed to be the dominant ions in sporadic E layers. Their chemistry, however, is not well known. For this reason and because large amounts of additional computer memory and time would be required for their inclusion in the transport mode, they have been omitted from the calculations presented here. This omission might explain why our electron density is somewhat low in the 90 to 110 km

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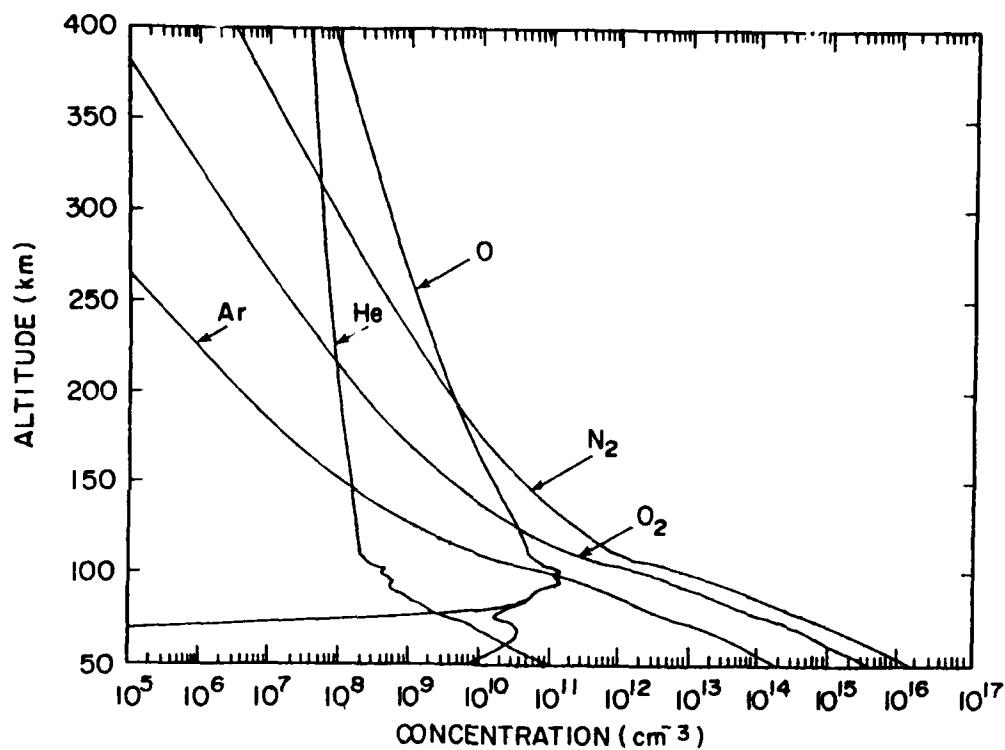


Figure 3. The Time Invariant Height Profiles of N_2 , O_2 , Helium, and Argon. Above about 100 km, atomic oxygen is also invariant. Below 100 km, the noon and midnight profiles of atomic oxygen are shown

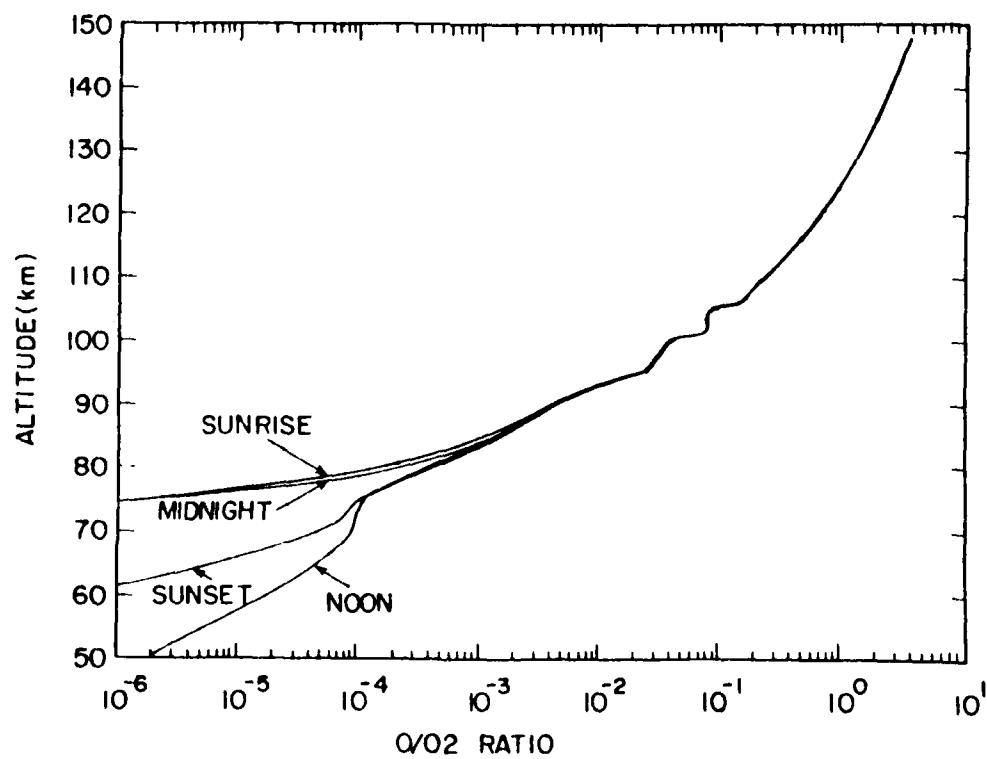


Figure 4. The Ratio of Atomic to Molecular Oxygen at Four Different Times During the Day. These curves reflect the diurnal variations of atomic oxygen in the mesosphere

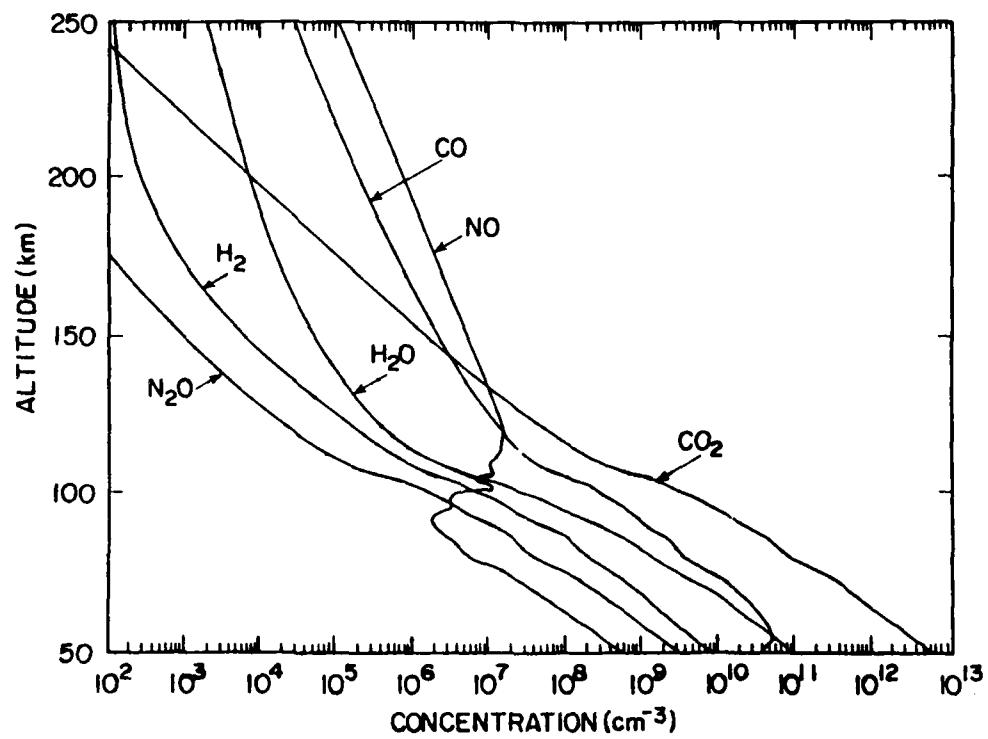


Figure 5. The Noontime Profiles of H_2 , H_2O , NO, N_2O , CO, and CO_2 in the Mesosphere and Lower Thermosphere

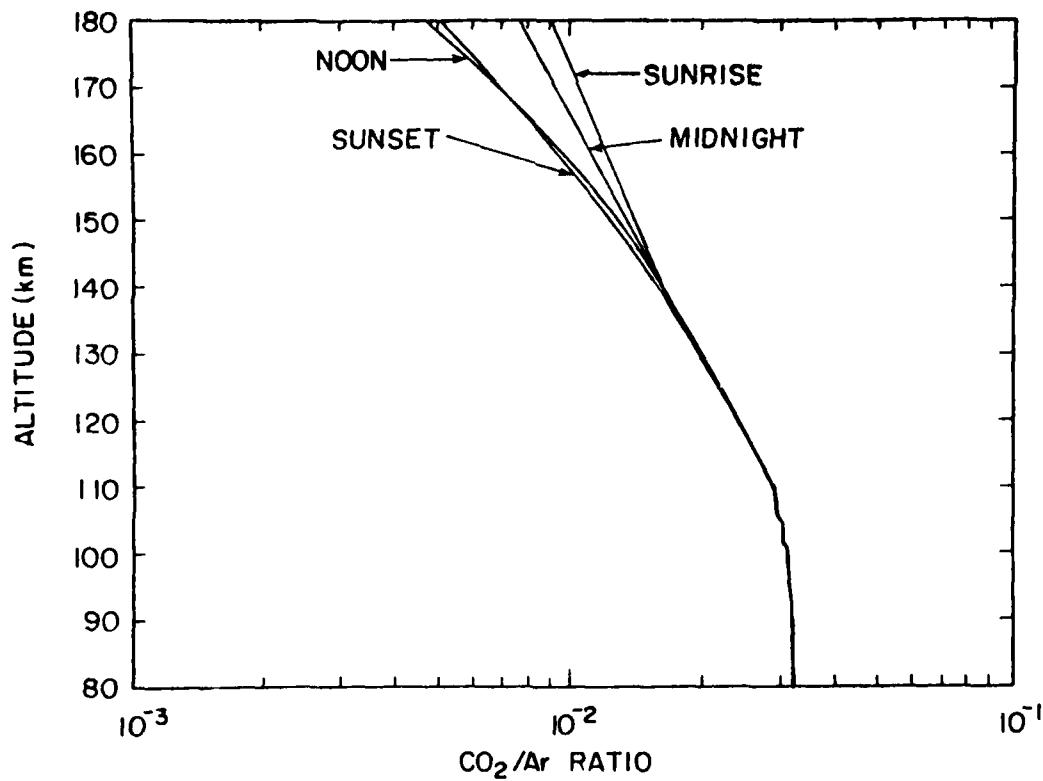


Figure 6. The Ratio of Carbon Dioxide to Argon at Four Different Times During the Day. The curves show the diurnal variation of carbon dioxide in the lower thermosphere

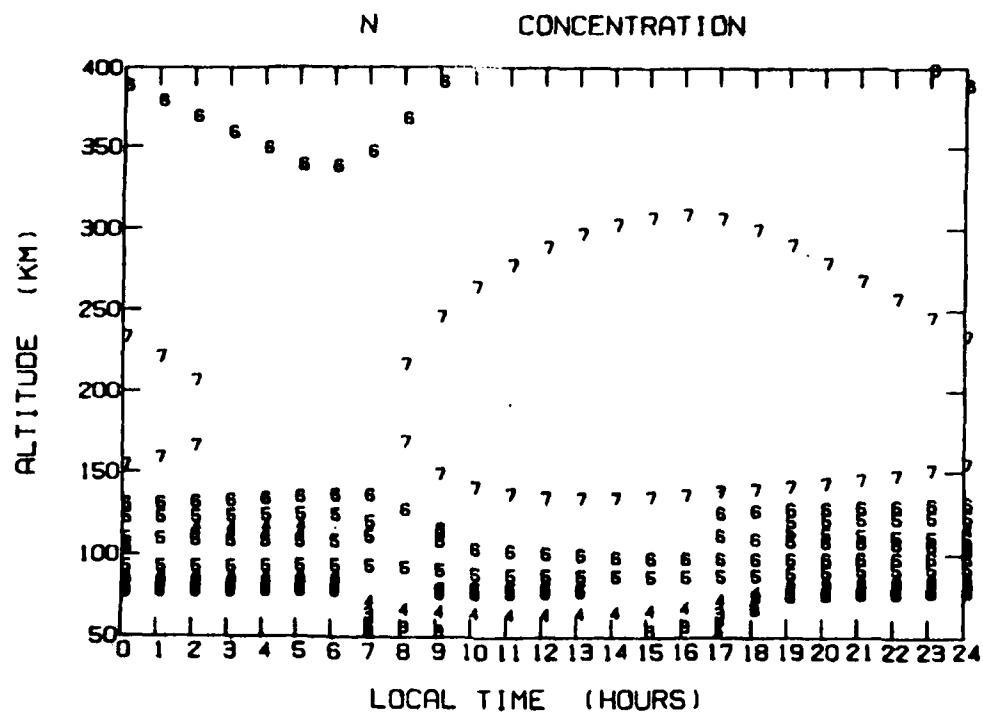


Figure 7. The Iso-Density Contours for Atomic Nitrogen

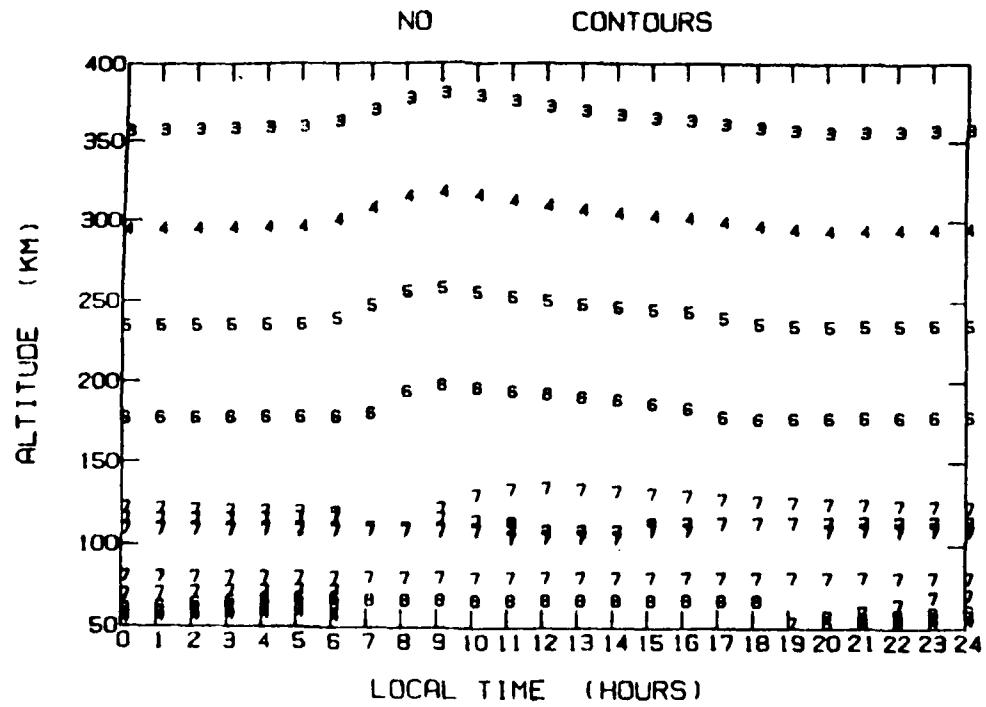


Figure 8. The Iso-Density Contours for Nitric Oxide

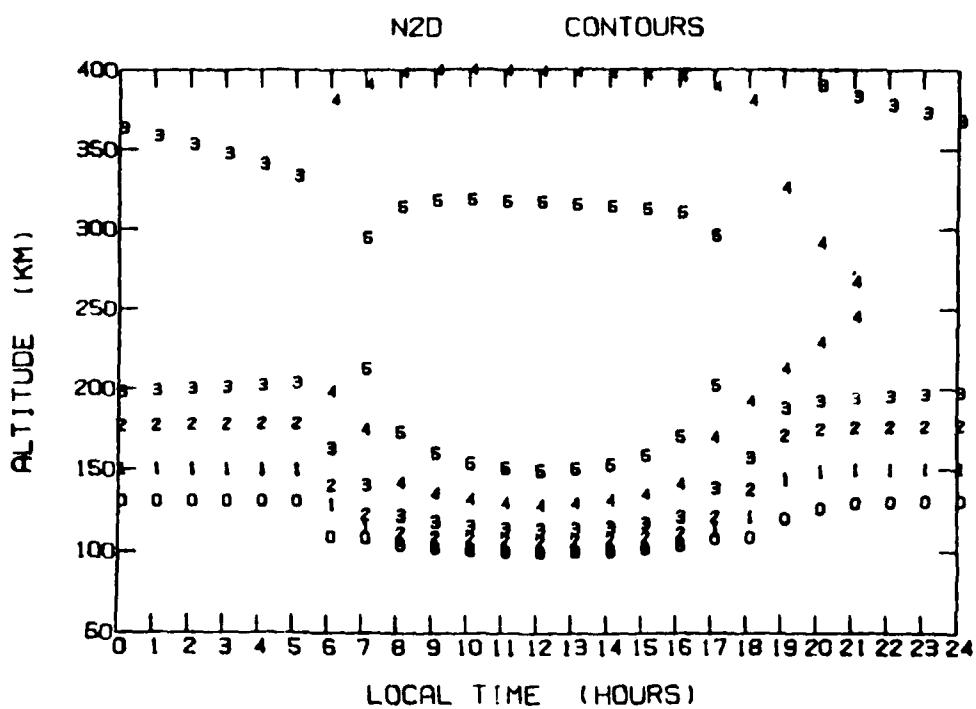


Figure 9. The Iso-Density Contours for N²D

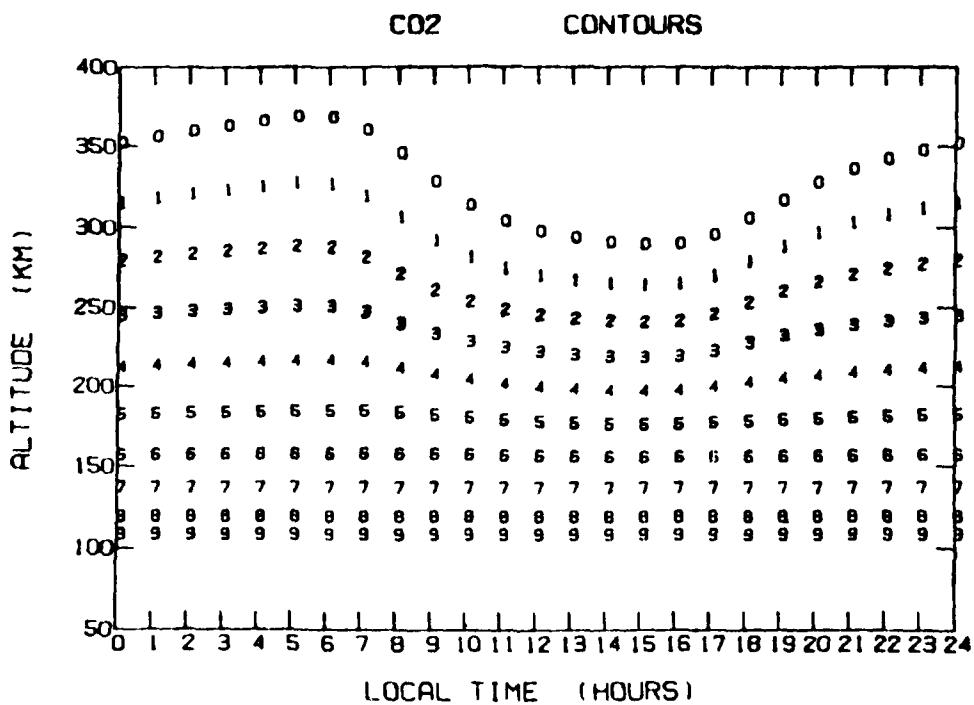


Figure 10. The Iso-Density Contours for Carbon Dioxide. The contours indicate a lack of any diurnal variation for this species below 150 km

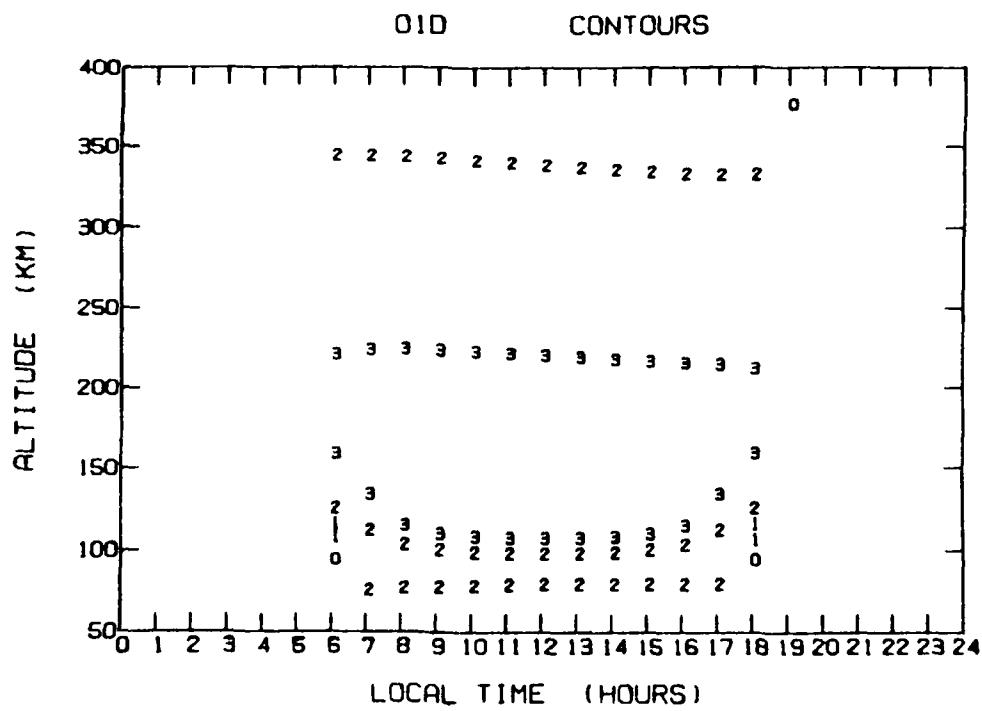


Figure 11. The Iso-Density Contours for O¹D

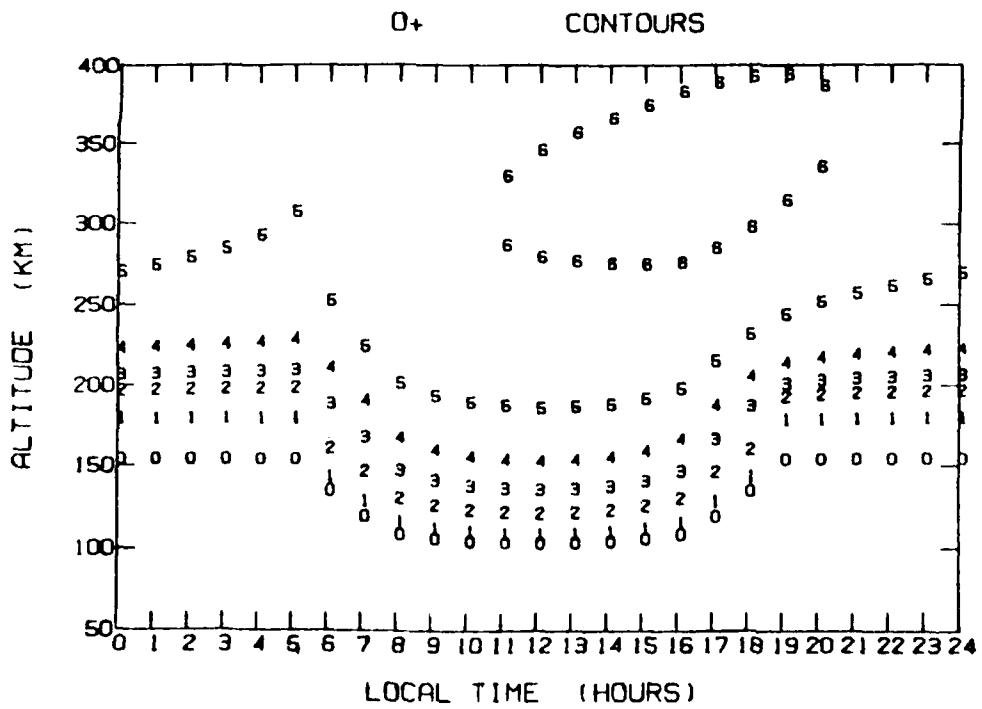


Figure 12. The Iso-Density Contours for O⁺. The contours show the daytime F-layer peak ionization in excess of 10^6 cm^{-3}

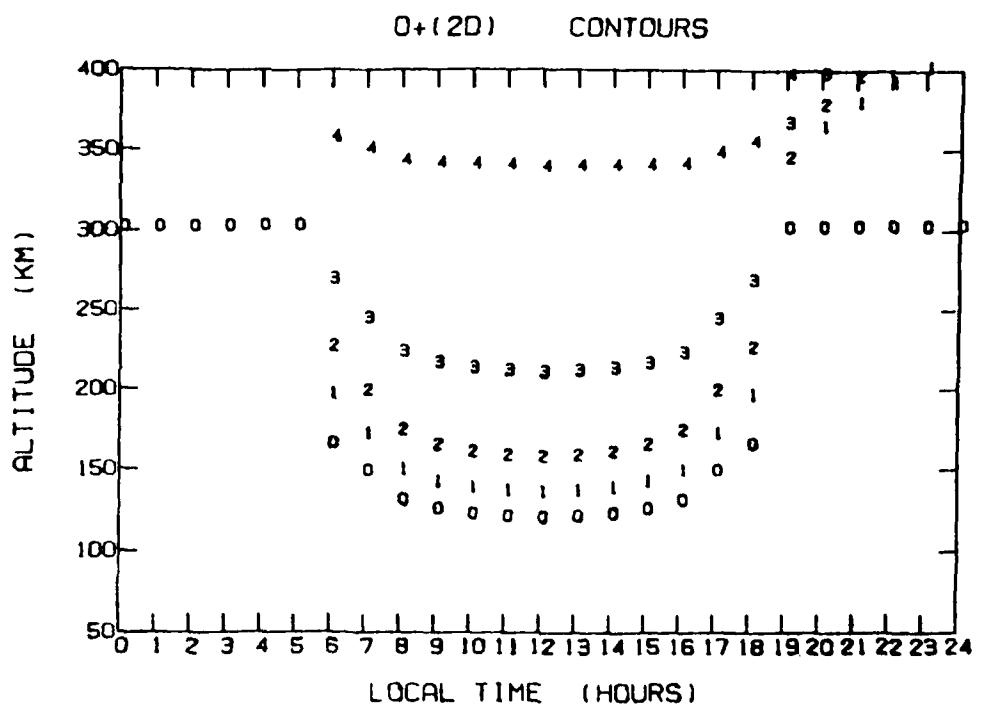


Figure 13. The Iso-Density Contours for $O^+(2D)$

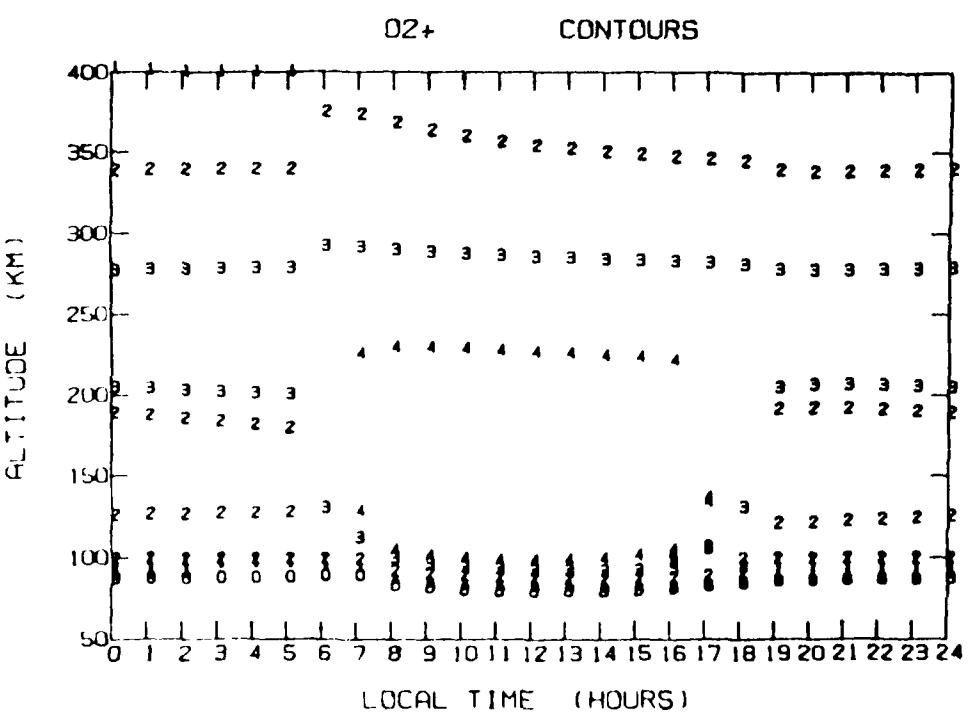


Figure 14. The Iso-Density Contours for O_2^+

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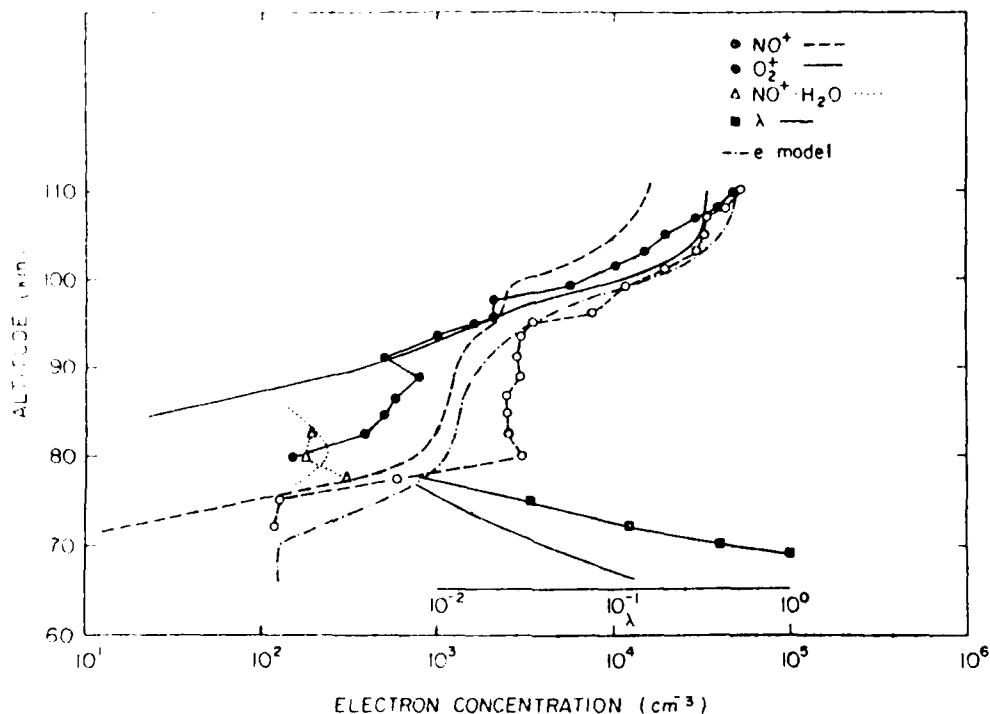
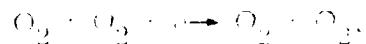


Figure 35. Altitude Profiles of NO_2^+ , O_2^+ , $\text{NO}_2^+\cdot\text{H}_2\text{O}$, Electrons and λ , the Ratio of Negative Ions to Electrons. The curves with symbols are the measurements reported by Torkar and Freidrich.⁴⁰ The curves without symbols are from this model

ions than the distribution determined in the Winter Anomaly campaign. However, there is a key reaction in the production of O_2^- in the early part of the negative ion chemistry scheme used by Viggiano et al.,⁴³ namely



They used a rate coefficient of about $1 \times 10^{-29} \text{ cm}^{-6} \text{ sec}^{-1}$ for this reaction. Phelps,^{A10} however, cites a rate coefficient of $1.4 \times 10^{-29} (300/T) \exp(-600/T)$ for this reaction. This value would result in a theoretical altitude distribution of the negative ions similar to that in our calculations. Therefore, these data of Arnold et al.⁴¹ still are unexplained by theory and leave us with D-region negative ions that are still largely not understood.

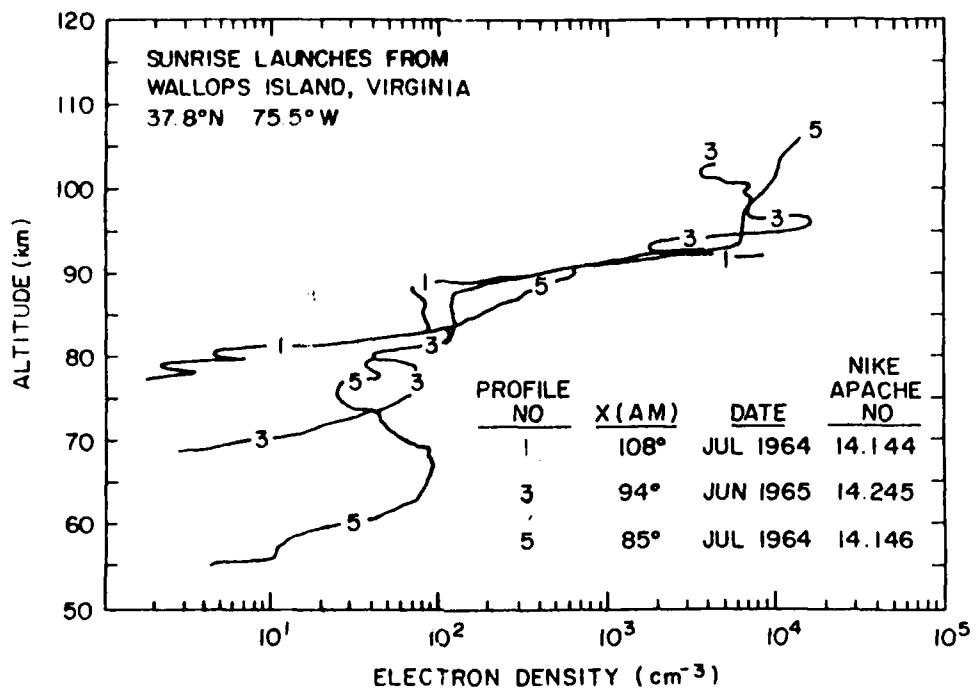


Figure 33. Altitude Profiles of Electron Density From Three Rocket Experiments at Wallops Island by Mechtly and Smith³⁹

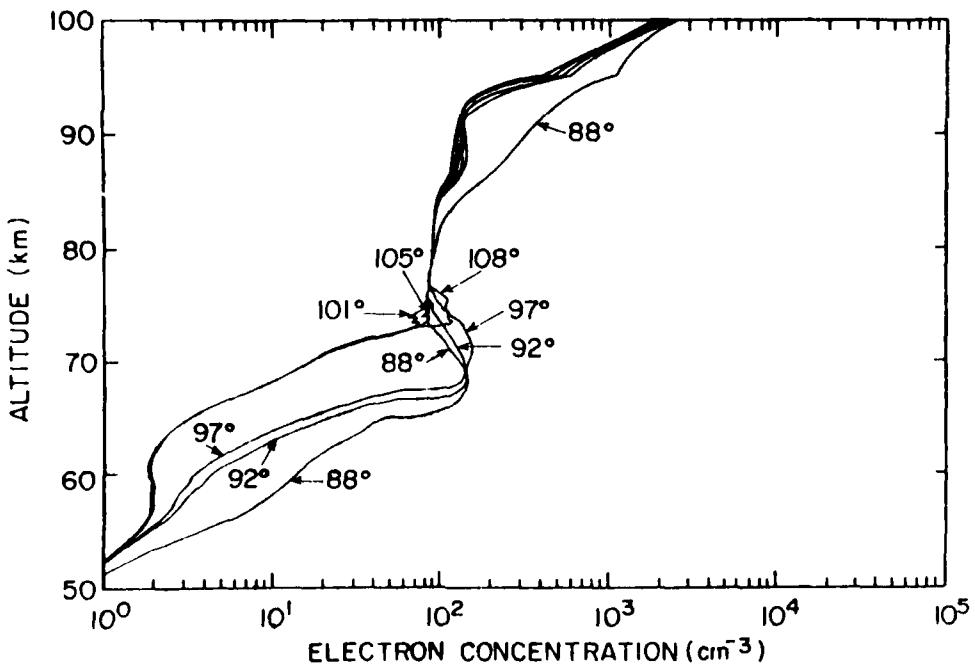


Figure 34. Altitude Profiles of the Electron Concentration From This Model at Several Zenith Angles During Sunrise

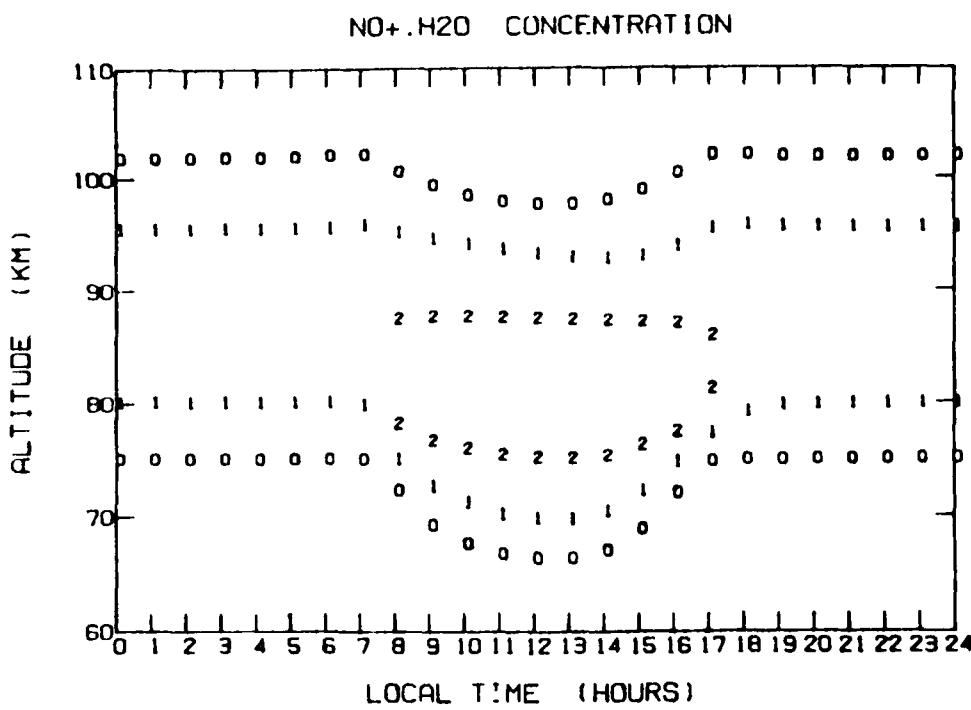


Figure 31. The Iso-Density Contours for NO⁺.H₂O in the Mesosphere

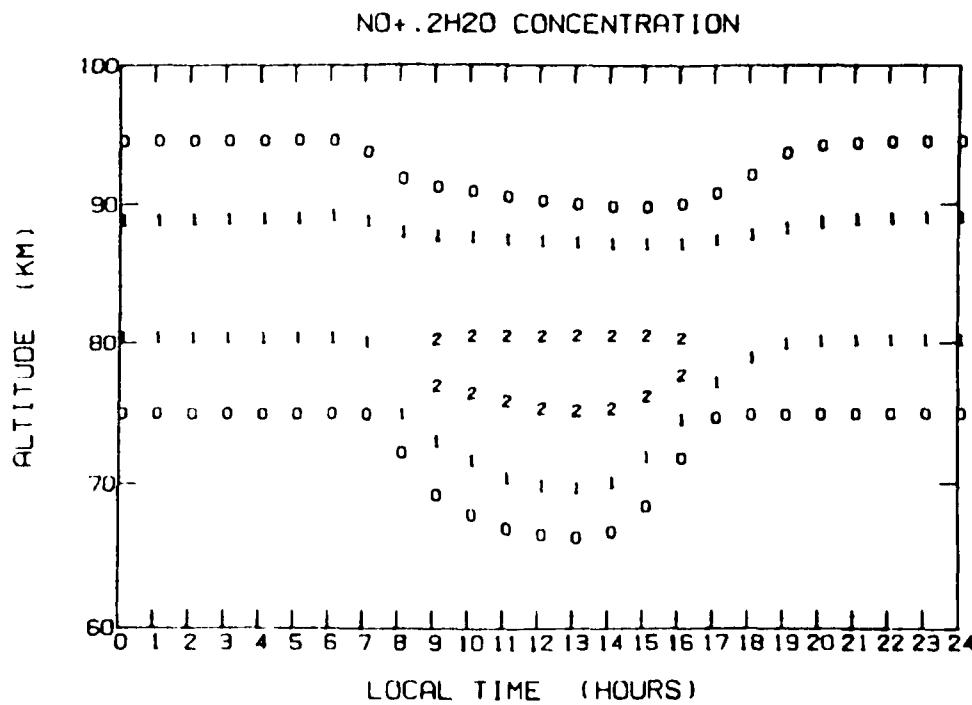


Figure 32. The Iso-Density Contours for NO⁺·(H₂O)₂ in the Mesosphere

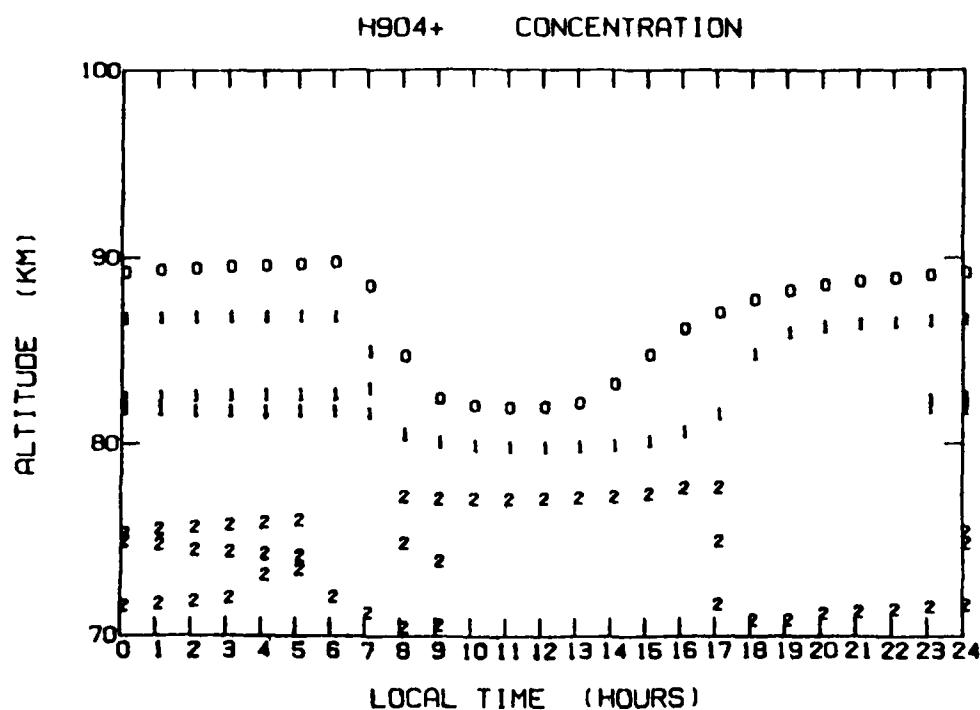


Figure 29. The Iso-Density Contours for H_9O_4^+ in the Mesosphere

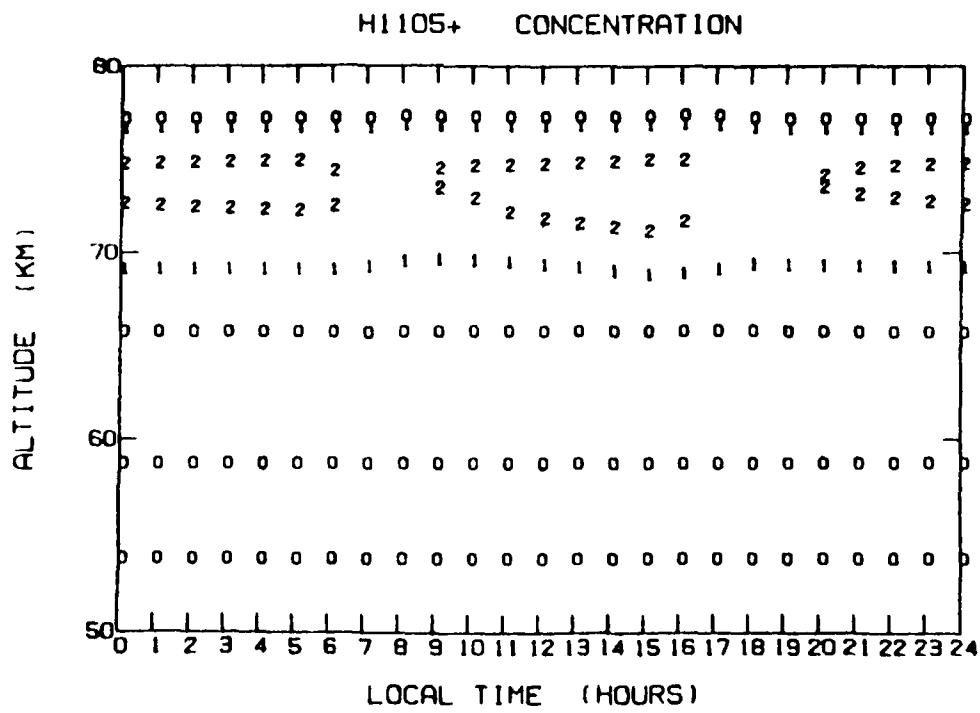


Figure 30. The Iso-Density Contours for $\text{H}_{11}\text{O}_5^+$ in the Mesosphere

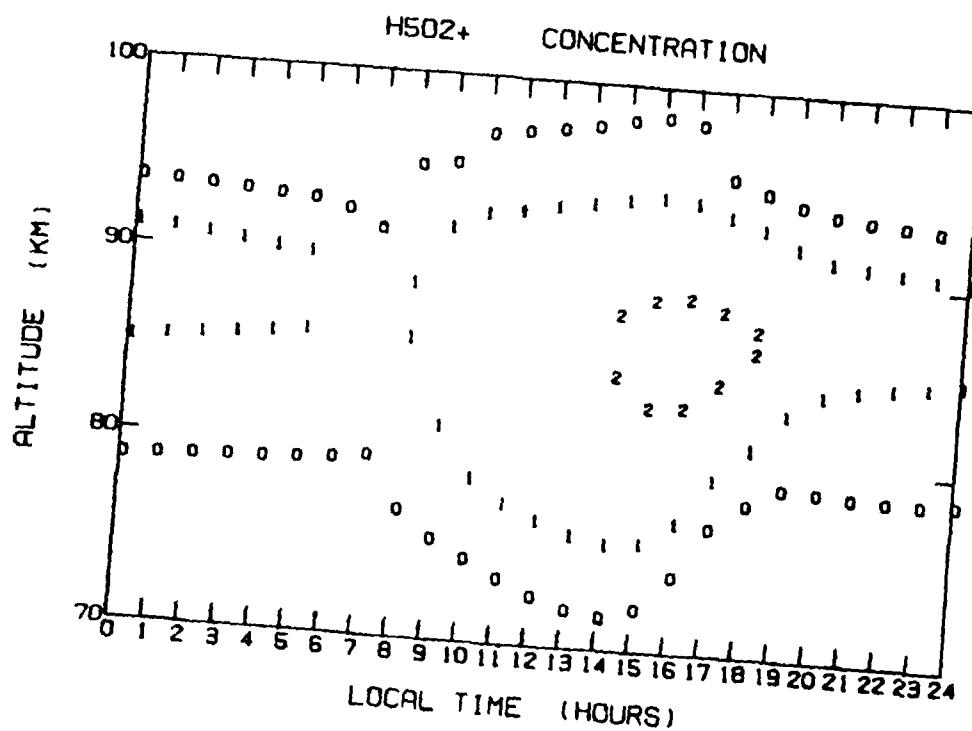


Figure 27. The Iso-Density Contours for H₅O₂⁺ in the Mesosphere

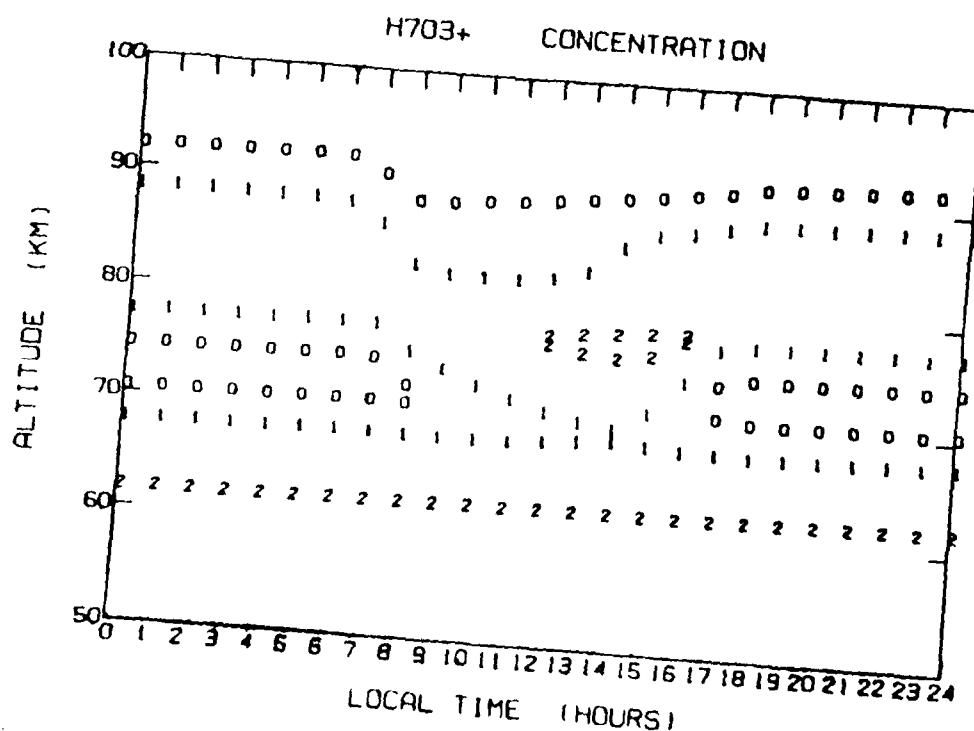


Figure 28. The Iso-Density Contours for H₇O₃⁺ in the Mesosphere

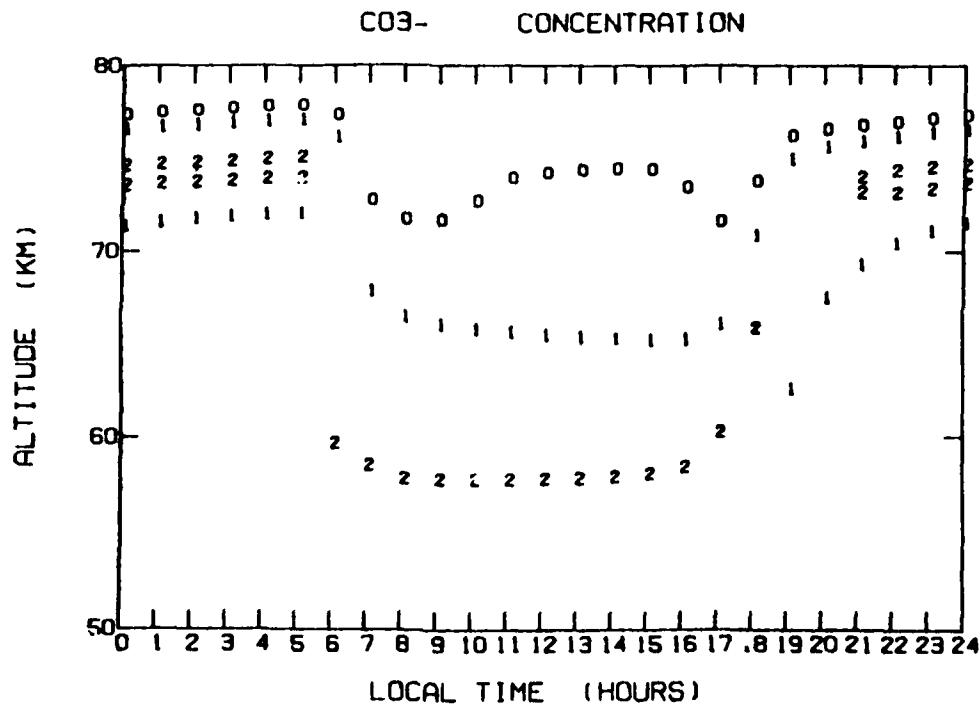


Figure 25. The Iso-Density Contours for CO₃⁻ in the Mesosphere

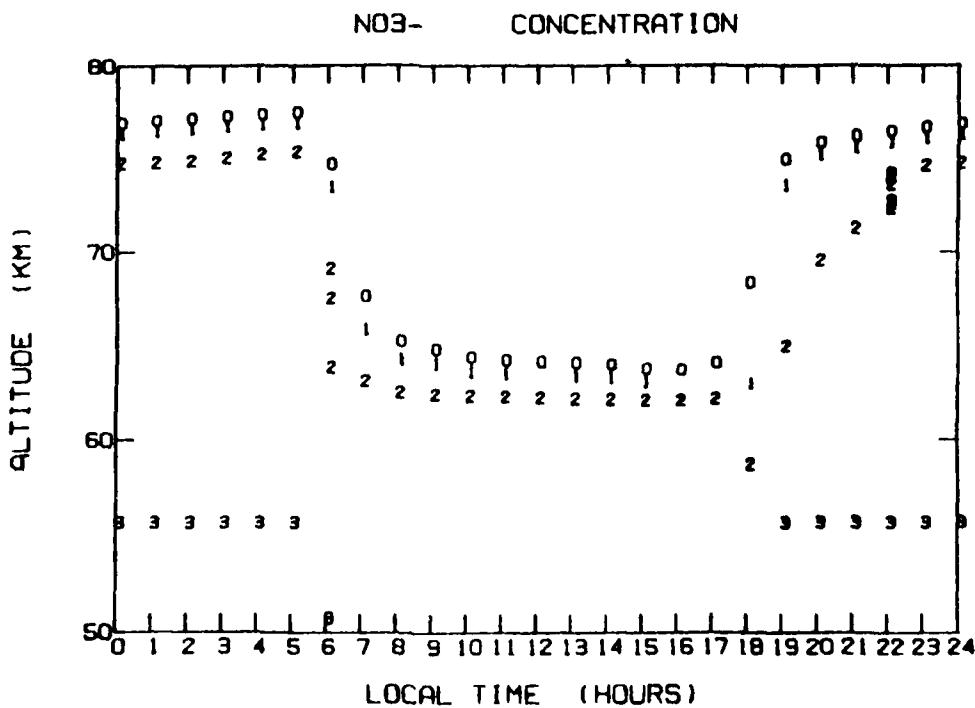


Figure 26. The Iso-Density Contours for NO₃⁻ in the Mesosphere

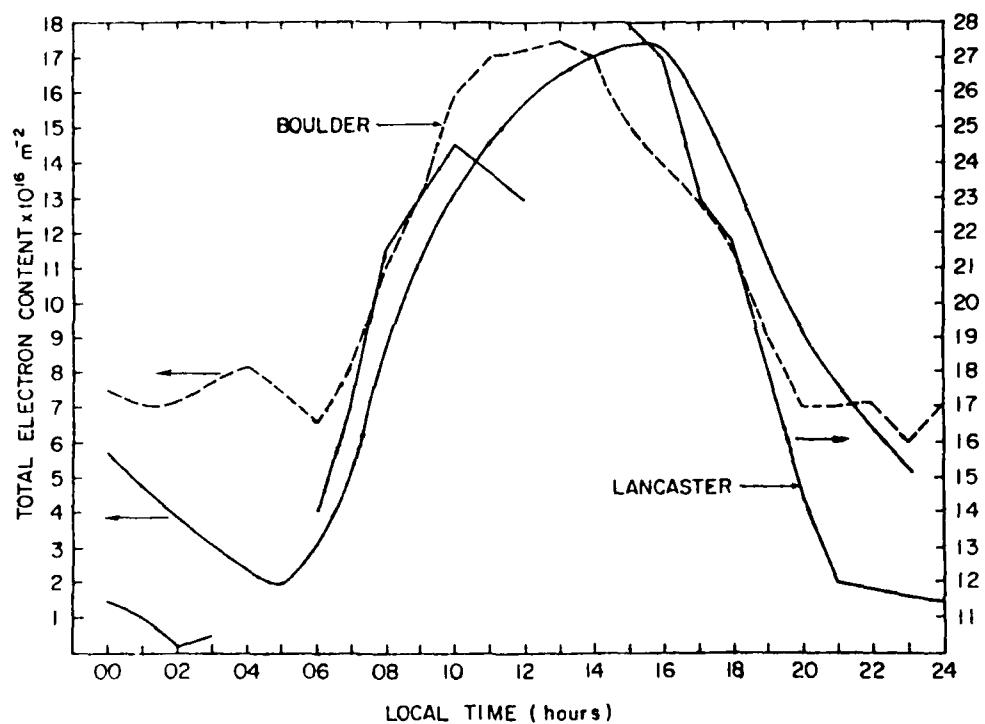


Figure 23. The Total Electron Content. The continuous curve is for this model. The dashed and broken curves are the measurements of Pouiter et al.³⁴

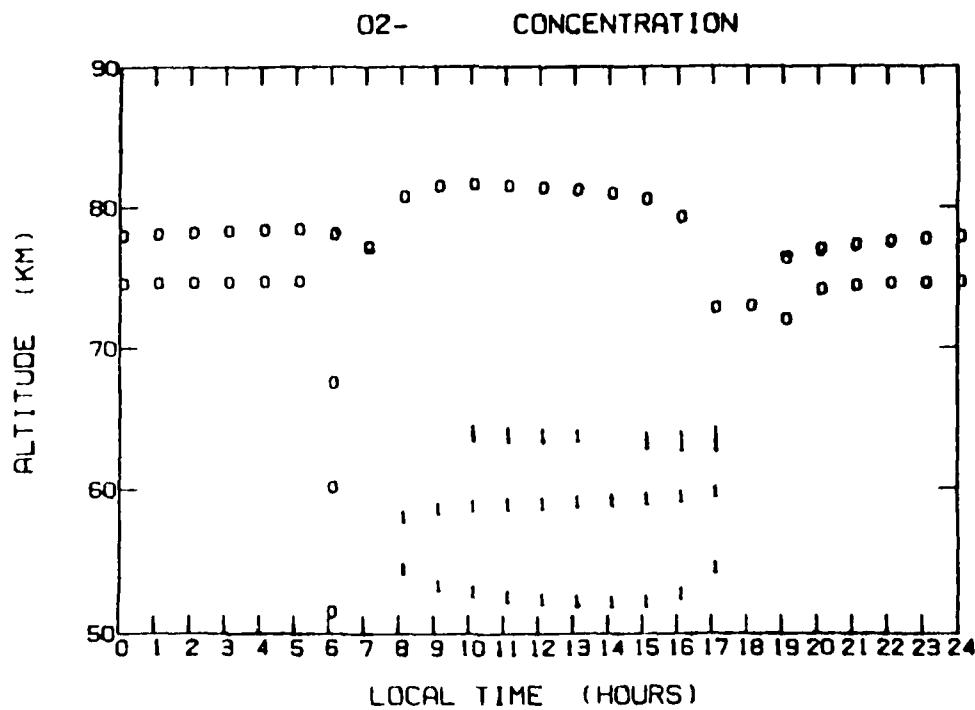


Figure 24. The Iso-Density Contours for O₂⁻ in the Mesosphere

region (Figure 19) compared to the work of Solomon et al.³⁸ Below 90 km, the cluster and negative ions change the effective electron density to a point where it disappears completely between 60 and 70 km.

The isodensity contours for the negative ions and the positive water cluster ions for the chemistry given in Appendix A are displayed in Figures 24 to 32. Iso-density plots for the remaining positive and negative mesospheric ions are not presented here because these ions never achieve a concentration greater than 1 cm^{-3} at any time during the diurnal cycle. Of course, the dominant negative ions in the mesosphere are apparently the cluster ions. Because their chemistry is still not fully understood, they have not been included in this work.

Particular emphasis is given to the sunrise period and a comparison of our electron density profiles with sequential measurements made at Wallops Island by Mechtly and Smith.³⁹ For clarity, the measured profiles are displayed in Figure 33, and our calculated profiles are shown in Figure 34. In some instances, the comparison of theory with experiment is fairly good. Substantial differences, however, do exist. These are, in part, a result of the differences in the seasons (summer for the measurement and winter for the model).

In explaining the D and lower E regions of the ionosphere, there is a question, still not completely answered, of the distribution of the positive and negative ions. To exemplify the puzzle, we cite two works. First, there is the report by Torkar and Freidrich⁴⁰ of the measurements taken during the "Winter Anomaly" campaign in Spain. The comparison between our calculations and those measurements, shown in Figure 35, is surprisingly good. In contrast are the measurements of Arnold et al⁴¹ and Viggiano and Arnold.⁴² These measurements, supposedly explained by Viggiano et al,⁴³ show an extremely different distribution of the negative

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38. Solomon, S., Reid, G.C., Roble, R.G., and Crutzen, P.J. (1982) Photochemical coupling between the thermosphere and the lower atmosphere, 2. D region ion chemistry and the winter anomaly, *J. Geophys. Res.* 87:7221-7227.
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 40. Torkar, K.M., and Friedrich, M. (1983) Tests of an ion-chemical model of the D- and lower E-region, *J. Atmos. Terr. Phys.* 45:369-385.
 41. Arnold, F., Viggiano, A.A., and Ferguson, E.E. (1982) Combined mass spectrometer composition measurements of positive and negative ions in the lower ionosphere-II. Negative ions, *Planet. Space Sci.* 30:1307-1314.
 42. Viggiano, A.A., and Arnold, F. (1981) The first height measurements of the negative ion composition of the stratosphere, *Planet. Space Sci.* 29:895-906.
 43. Viggiano, A.A., Arnold, F., Fahey, D.W., Fehsenfeld, F.C., and Ferguson, E.E. (1982) Silicon negative ion chemistry in the atmosphere--in situ and laboratory measurements, *Planet. Space Sci.* 30:499-506.

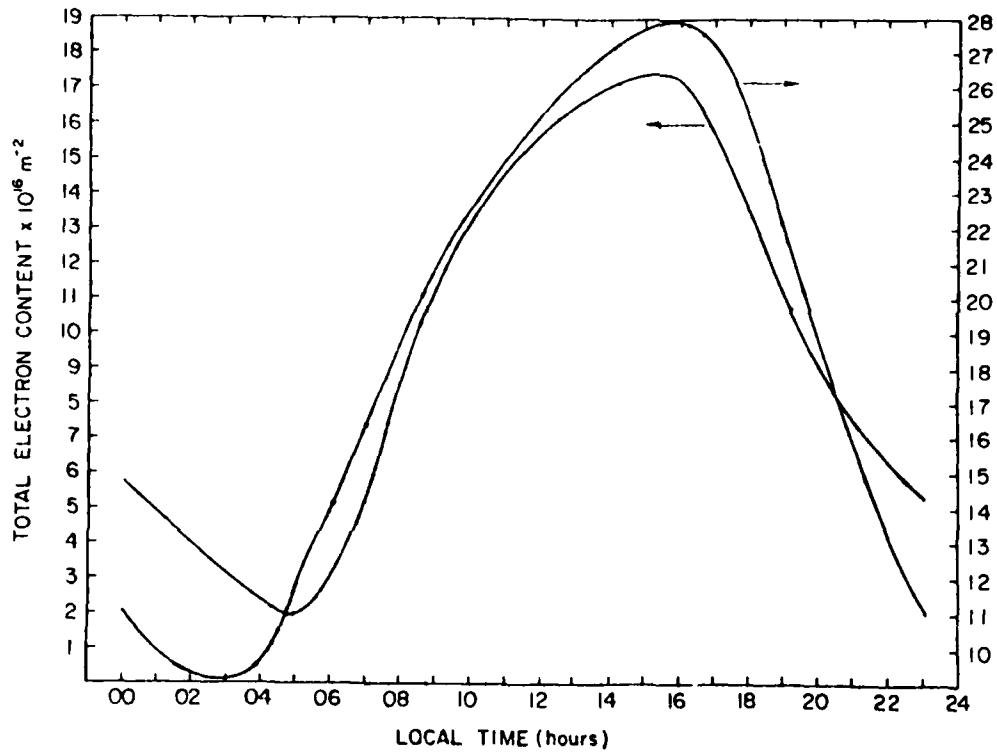


Figure 21. The Total Electron Content. The left-hand scale is for this model. The right-hand scale is for the calculations of Sethia et al³³

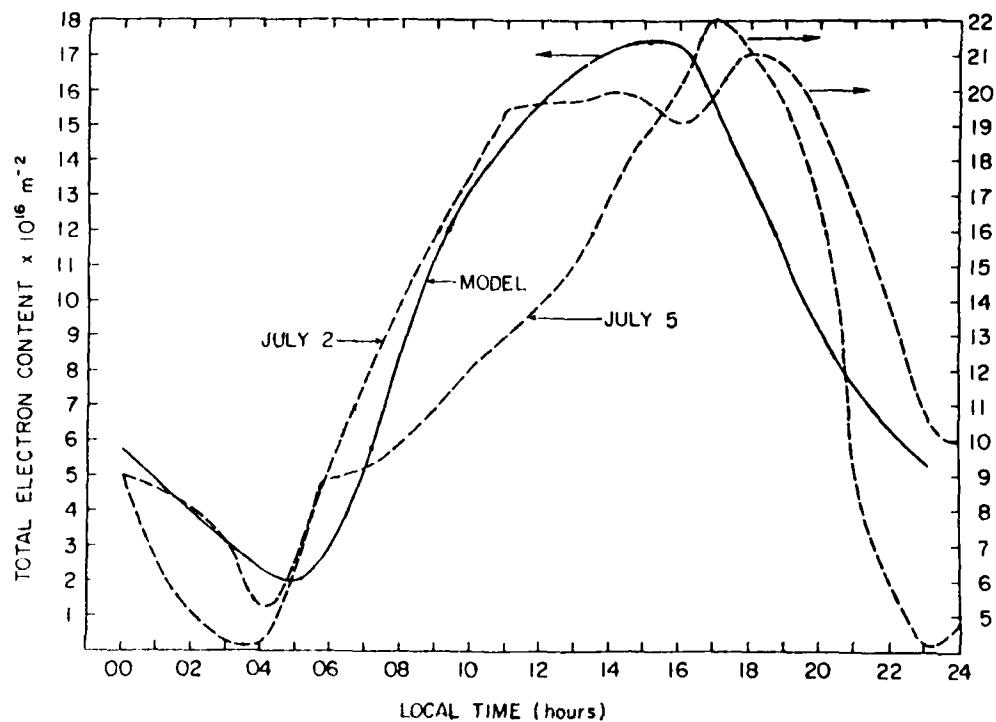


Figure 22. The Total Electron Content. The solid curve is for this model. The dashed curves are the measurements of Sethia et al³³

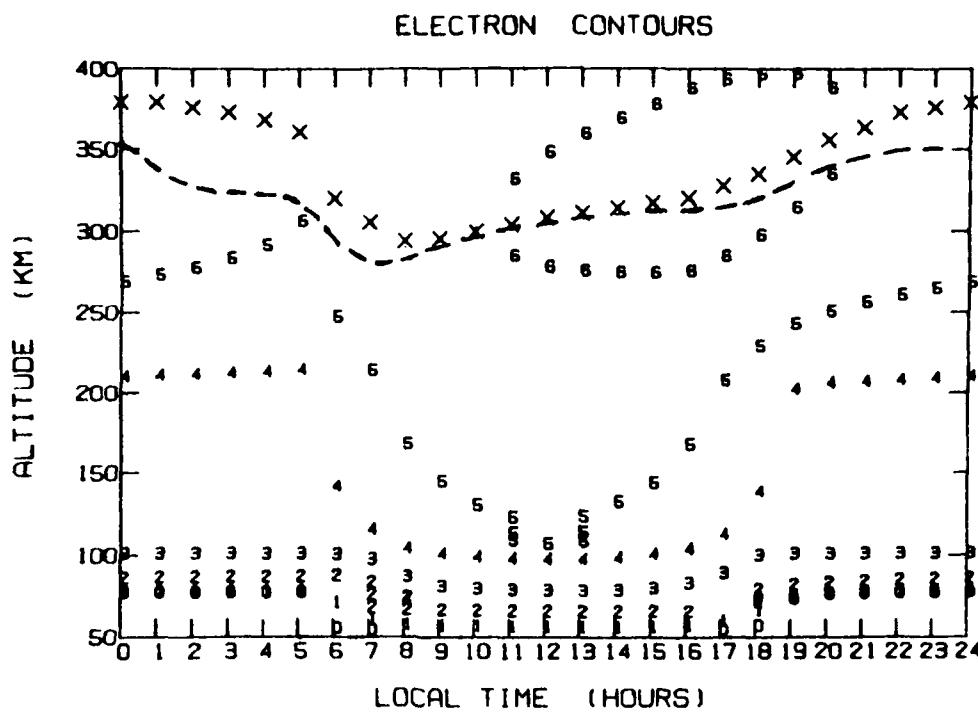


Figure 19. The Iso-Density Contours for the Electrons. The crosses show the diurnal change in the altitude of the peak electron concentration in the model. The dashed curve indicates the measurement of this parameter by Evans³¹

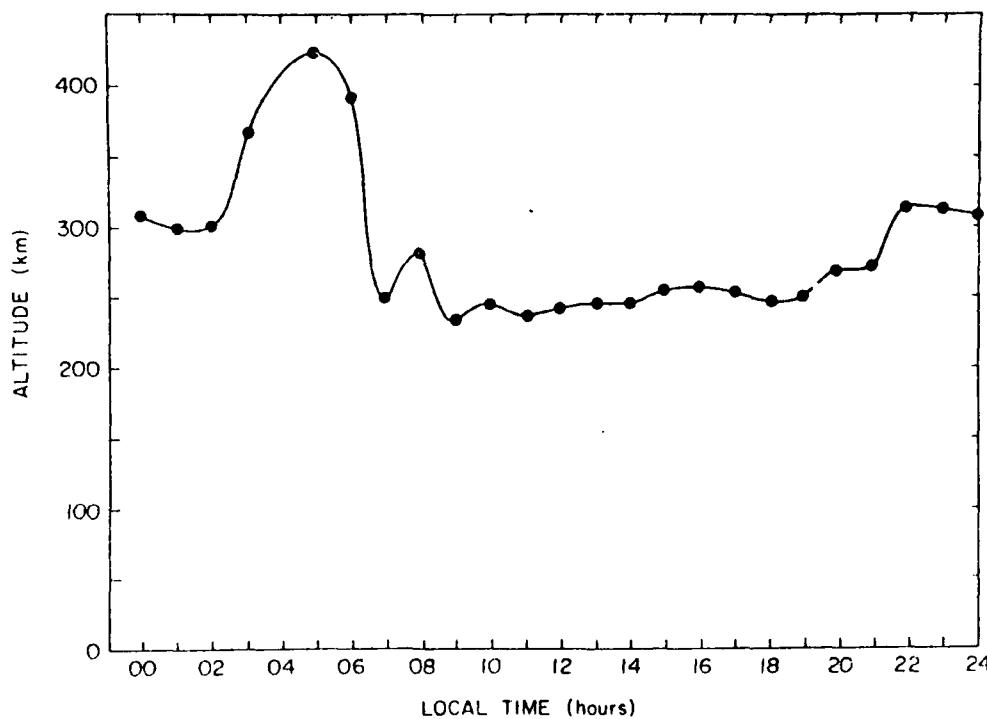


Figure 20. Measurements of Reinisch³² of the Height of the F-Layer Peak at Eglin AFB, Fla., on 20/21 November 1970

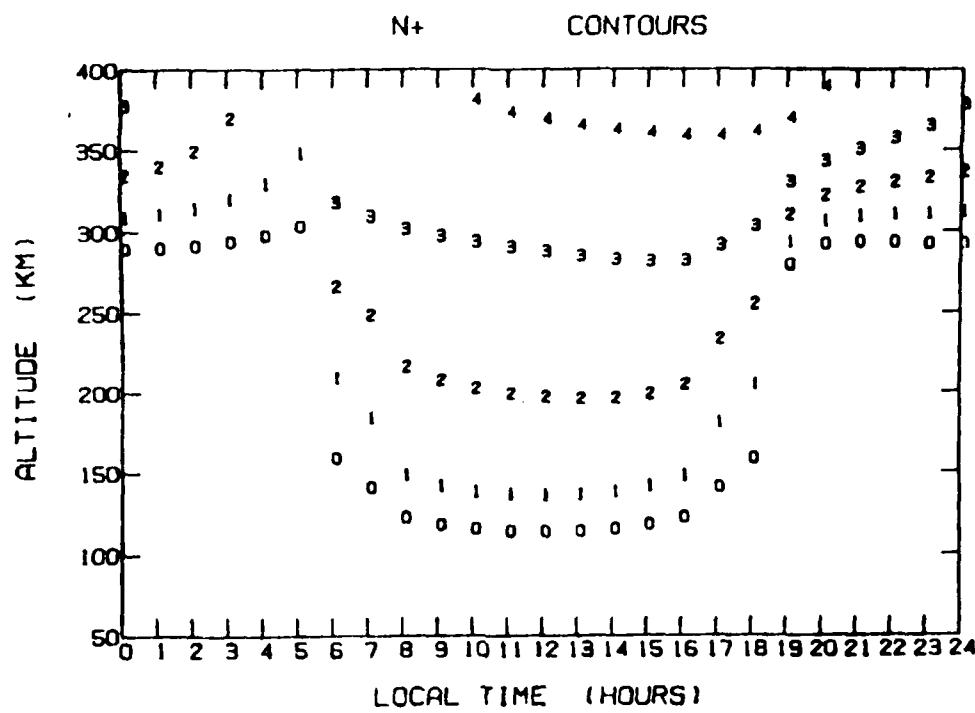


Figure 17. The Iso-Density Contours for N^+

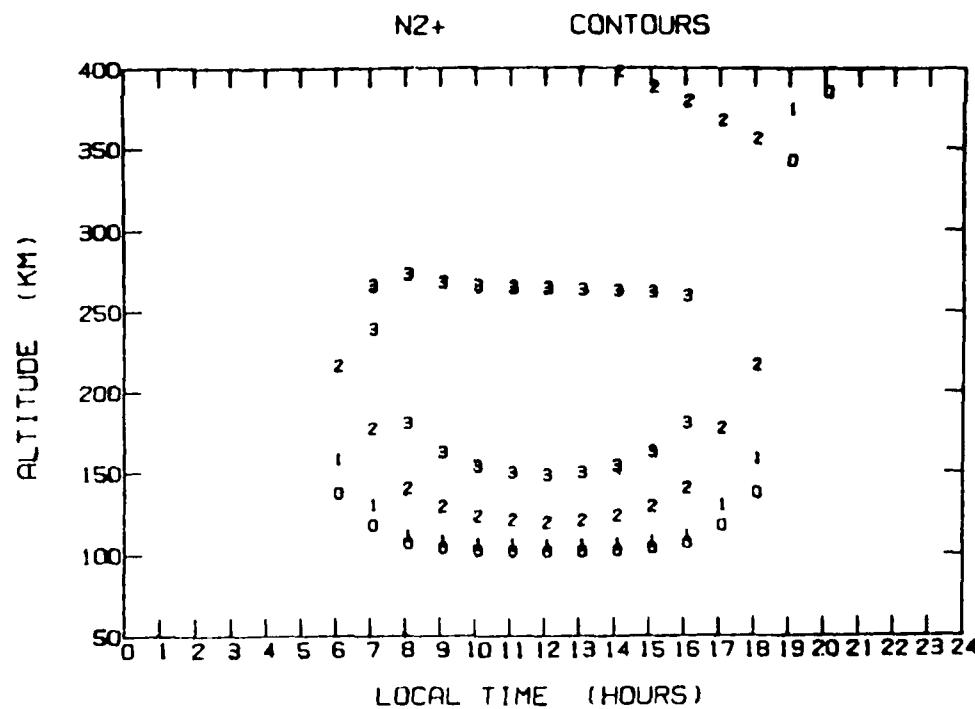


Figure 18. The Iso-Density Contours for N_2^+ . At night, the concentration of N_2^+ falls below 1 cm^{-3} at all altitudes

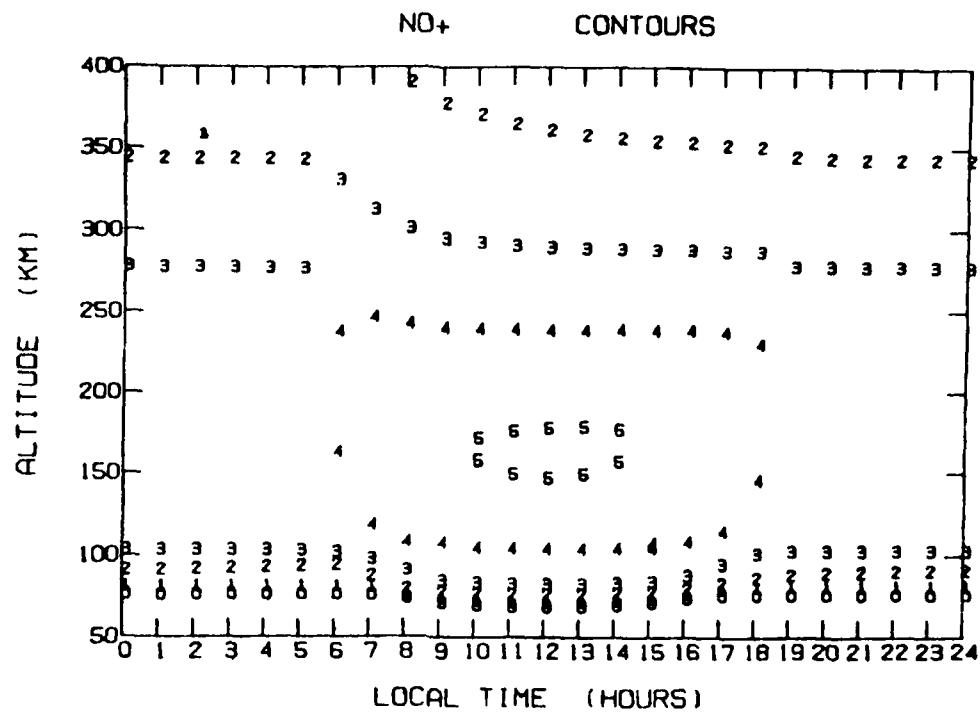


Figure 15. The Iso-Density Contours for NO⁺

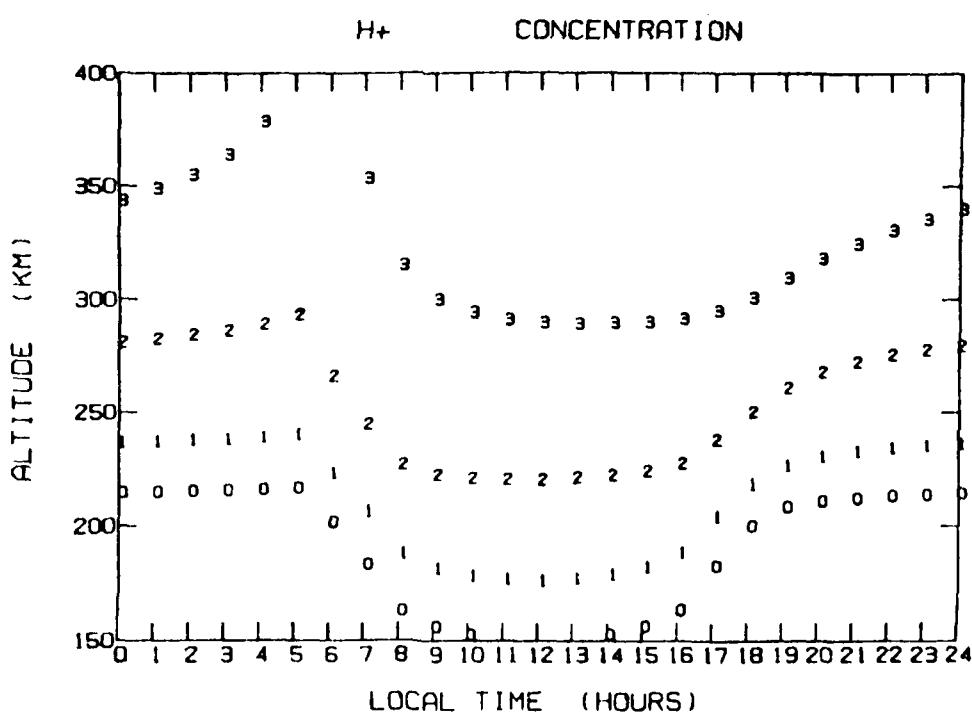


Figure 16. The Iso-Density Contours for H⁺

34. Poulter, E.M., Hargreaves, J.K., Bailey, G.J., and Moffet, R.J. (1981) Electron content modelling: The significance of protonospheric contents, Planet. Space Sci. 29:869-883.
35. Kohl, H., King, J.W., and Eccles, D. (1969) An explanation of the magnetic declination effect in the ionospheric F-layer, J. Atmos. Terr. Phys. 31:1011-1016.
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38. Solomon, S., Reid, G.C., Roble, R.G., and Crutzen, P.J. (1982) Photochemical coupling between the thermosphere and the lower atmosphere, 2. D region ion chemistry and the winter anomaly, J. Geophys. Res. 87:7221-7227.
39. Mechtly, E.A., and Smith, L.G. (1968) Growth of the D-region at sunrise, J. Atmos. Terr. Phys. 30:363-369.
40. Torkar, K.M., and Friedrich, M. (1983) Tests of an ion-chemical model of the D- and lower E-region, J. Atmos. Terr. Phys. 45:369-385.
41. Arnold, F., Viggiano, A.A., and Ferguson, E.E. (1982) Combined mass spectrometer composition measurements of positive and negative ions in the lower ionosphere—II. Negative ions, Planet. Space Sci. 30:1307-1314.
42. Viggiano, A.A., and Arnold, F. (1981) The first height measurements of the negative ion composition of the stratosphere, Planet. Space Sci. 29:895-906.
43. Viggiano, A.A., Arnold, F., Fahey, D.W., Fehsenfeld, F.C., and Ferguson, E.E. (1982) Silicon negative ion chemistry in the atmosphere—in situ and laboratory measurements, Planet. Space Sci. 30:499-506.

Appendix A

The Chemical Reactions

Table A1 lists the complete set of chemical reactions used in this report. Reactions 1 through 127 have temperature invariant rate coefficients that are listed after each reaction along with the source from which the rate coefficient was taken. Reactions 128 through 184 are temperature dependent. The rate coefficients for these reactions are of the form

$$k(z) = A \left(\frac{T(z)}{300} \right)^B \exp \left(\frac{C}{T(z)} \right) \quad (A1)$$

where the A's, B's, and C's are listed in that order after each reaction and T(z) is the temperature at altitude z. Reactions 185 through 207 are the photo-reactions, and are functions of the wavelength of the solar radiation and the local solar flux. Reactions 208 through 215 are the energetic electron processes. Some of the rate coefficients in the list of reactions have not been measured in the laboratory. Therefore, these rate coefficients are given as estimates.

Table A1. Chemical Reaction Scheme

NO.	CHEMICAL REACTION	REFERENCE			
		A	B	C	
1	O + N2D	+ O	1.00E-12	0.00	
2	O + OH	+ H	2.20E-11	0.00	
3	O + HO2	+ OH	1.00E-11	0.00	
4	O + NO2	+ NO	9.10E-12	0.00	
5	H + HO2	+ H2	1.40E-11	0.00	
6	H + HO2	+ OH	1.00E-11	0.00	
7	H + O3	+ OH	2.60E-11	0.00	
8	N + O3	+ O2	5.70E-13	0.00	
9	N + NO2	+ N2O	8.00E-12	0.00	
10	N + NO2	+ NO	6.00E-12	0.00	
11	N + NO2	+ N2	4.00E-12	0.00	
12	CH + OH	+ H2O	2.00E-12	0.00	
13	OH + HO2	+ H2O	2.00E-10	0.00	
14	N2D + O2	+ OH	6.00E-12	0.00	
15	O1D + H2O	+ OH	2.10E-10	0.00	
16	O1D + H	+ OH	3.00E-11	0.00	
17	O1D + H2	+ H	1.30E-10	0.00	
18	O2D + H	+ H	4.40E-19	0.00	
19	O2D + H	+ H2	2.58E-04	0.00	
20	HNO2 + O	+ OH	1.00E-15	0.00	
21	HNO2 + OH	+ H2O	6.60E-12	0.00	
22	N2 + E	+ N	1.50E-07	0.00	
23	N2 + E	+ N2D	1.50E-07	0.00	
24	H3O+	+ H2	1.00E-06	0.00	
25	H5O2+	+ E	2.00E-06	0.00	
26	H7O3+	+ E	4.00E-06	0.00	
27	H9O4+	+ E	5.00E-06	0.00	
28	H11O5+	+ E	7.50E-06	0.00	
29	H3O+·OH	+ E	2.00E-06	0.00	
30	H5O2+·CO2+	+ E	2.00E-06	0.00	
31	O + E	O-	1.30E-15	0.00	
32	O2- + N2	+ N2	1.00E-31	0.00	
33	O2 + NO	+ O2	6.00E-10	0.00	
34	O2 + N	+ NO+	1.80E-10	0.00	
35	O2 + NO2	+ O2	6.60E-10	0.00	
36	O4 + O2	+ O2+	3.00E-10	0.00	
37	O4 + H2O	+ O2+	1.50E-09	0.00	
38	N2 + O	+ NO+	7.50E-11	0.00	
39	N2 + O	+ NO+	6.00E-10	0.00	
40	O + H	+ O	4.00E-10	0.00	
41	H + O	+ H	6.00E-10	0.00	
42	N + O2	+ ND+	6.00E-11	0.00	
43	N + O2	+ N	6.00E-10	0.00	
44	N + O2	+ C2+	2.00E-10	0.00	
45	HE + O2	+ O	+ HE	1.50E-09	0.00
46	HE + O2	+ C2+	+ HE	2.00E-10	0.00
47	HE + N2	+ N	+ HE	1.50E-09	0.00

Table A1. Chemical Reaction Scheme (Contd)

NO.	CHEMICAL REACTION		A	B	C	REFERENCE
	Reactants	Products				
48	HE + N2	+ HE	3.50E-10	0.00	0.	A24 FERGUSON (1974)
49	HE + NO	+ N+	1.50E-09	0.00	0.	A26 FEHSENFIELD ET AL (1966)
50	NO2 + NO	+ O2	2.90E-10	0.00	0.	A26 FEHSENFIELD ET AL (1969A)
51	H5O2+ + H2O	+ M	3.00E-30	0.00	0.	ESTIMATED
52	H9O4+ + H2O	+ M	1.00E-27	0.00	0.	A23 HOWARD ET AL (1972)
53	O2 + H2O	+ H2O	1.00E-09	0.00	0.	A23 HOWARD ET AL (1972)
54	O2 + H2O	+ H2O	2.00E-10	0.00	0.	A26 PUCKETT AND TEAGUE (1971)
55	NO + 3H2O	+ H2O	7.00E-11	0.00	0.	A27 DUNKIN ET AL (1971)
56	NO + CO2	+ H2O	1.00E-09	0.00	0.	ESTIMATED
57	H3O+ . N2	+ CO2	1.00E-10	0.00	0.	ESTIMATED
58	H3O+ . CO2	+ H2O	1.00E-27	0.00	0.	ESTIMATED
59	H5O2+ . CO2	+ H2O	1.00E-10	0.00	0.	ESTIMATED
60	H3O+ . CO2	+ H2O	1.00E-10	0.00	0.	ESTIMATED
61	H3O+ . OH	+ H2O	1.40E-09	0.00	0.	A23 HOWARD ET AL (1972)
62	H3O+ + NO2-	+ H2O	6.00E-08	0.00	0.	ESTIMATED
63	H3O+ + NO3-	+ H2O	6.00E-08	0.00	0.	ESTIMATED
64	H3O+ + CO3-	+ H2O	6.00E-08	0.00	0.	ESTIMATED
65	H5O2+ + O2-	+ H2O	6.00E-08	0.00	0.	ESTIMATED
66	H5O2+ + NO2-	+ 2H2O	+ NO2	+ HO2	6.00E-08	0.00
67	H5O2+ + NO3-	+ 2H2O	+ NO2	+ OH	6.00E-08	0.00
68	H5O2+ + CO3-	+ 2H2O	+ NO2	+ OH	6.00E-08	0.00
69	H5O2+ + CO4-	+ 2H2O	+ CO2	+ HO2	6.00E-08	0.00
70	H7O3+ + NO2-	+ 3H2O	+ NO2	+ H	6.00E-08	0.00
71	H7O3+ + NO3-	+ 3H2O	+ NO2	+ OH	6.00E-08	0.00
72	H7O3+ + CO3-	+ 3H2O	+ CO2	+ OH	6.00E-08	0.00
73	H7O3+ + CO4-	+ 3H2O	+ CO2	+ HO2	6.00E-08	0.00
74	H9O4+ + NO2-	+ 4H2O	+ NO2	+ H	6.00E-08	0.00
75	H9O4+ + NO3-	+ 4H2O	+ NO2	+ OH	6.00E-08	0.00
76	H9O4+ + CO3-	+ 4H2O	+ CO2	+ OH	6.00E-08	0.00
77	H9O4+ + CO4-	+ 4H2O	+ CO2	+ HO2	6.00E-08	0.00
78	H11O5+ + NO2-	+ 5H2O	+ NO2	+ NO	6.00E-08	0.00
79	H11O5+ + NO3-	+ 5H2O	+ NO2	+ NO2	6.00E-08	0.00
80	H11O5+ + CO3-	+ 5H2O	+ CO2	+ OH	6.00E-08	0.00
81	H11O5+ + CO4-	+ 5H2O	+ CO2	+ HO2	6.00E-08	0.00
82	NO + H2O	+ H2O	+ NO2	+ C2	6.00E-08	0.00
83	NO + H2O	+ H2O	+ NO2	+ NO	6.00E-08	0.00
84	NO + H2O	+ CO3-	+ H2O	+ CO2	6.00E-08	0.00
85	NO + 2H2O	+ NO2-	+ 2H2O	+ NO2	6.00E-08	0.00
86	NO + 2H2O	+ NO3-	+ 2H2O	+ NO2	6.00E-08	0.00
87	NO + 2H2O	+ CO3-	+ 2H2O	+ NO2	6.00E-08	0.00
88	NO + 3H2O	+ NO2-	+ 3H2O	+ NO2	6.00E-08	0.00
89	NO + 3H2O	+ NO3-	+ 3H2O	+ NO2	6.00E-08	0.00
90	H9 + 3H2O	+ CO3-	+ 3H2O	+ C2	6.00E-08	0.00
91	H3O+ . CO2	+ NO3-	+ H2O2	+ NO2	6.00E-08	0.00
92	H3O+ . CO2	+ CO3-	+ CO2	+ OH	6.00E-08	0.00
93	H3O+ . CO2	+ CO4-	+ CO2	+ HO2	6.00E-08	0.00
94	H5O2+ . CO2	+ CO3-	+ 2CO2	+ OH	6.00E-08	0.00

Table A1. Chemical Reaction Scheme (Contd)

NO.	CHEMICAL REACTION	C			REFERENCE	
		A	B	C		
95	H ₂ O ₂ * + CO ₂ * + CO ₄ -	+ 2CO ₂	+ H ₂ O ₂	6.00E-08	0.00	A30 TRANSCOMB ET AL (1958)
96	O ⁻	+ HV	+ E	1.40E+00	0.00	A31 FEHSENFELD ET AL (1967)
97	O-	+ O	+ E	2.00E-10	0.00	A32 FEHSENFELD ET AL (1967)
98	O-	+ O ₃	+ O	5.30E-10	0.00	A33 WOO ET AL (1969)
99	O ₂ -	+ HV	+ O ₂	3.30E-01	0.00	A34 FEHSENFELD ET AL (1969C)
100	O ₂ -	+ O	+ E	1.50E-10	0.00	A35 BORTNER AND BAURER (1979)
101	O ₂ -	+ O	+ O ₂	1.50E-10	0.00	A36 FEHSENFELD ET AL (1967B)
102	O ₂ -	+ O ₃	+ O ₂	4.00E-10	0.00	A37 FEHSENFELD ET AL (1969C)
103	O ₂ -	+ O ₂ 10	+ O ₂	2.00E-10	0.00	A38 WONG ET AL (1972)
104	O ₃ -	+ HV	+ E	7.00E-02	0.00	A39 PETERSON (1976)
105	O ₃ -	+ HV	+ O ₂	5.00E-01	0.00	A40 BORTNER AND BAURER (1979)
106	O ₃ -	+ O	+ O ₂	3.20E-10	0.00	A41 FEHSENFELD ET AL (1969B)
107	O ₄ -	+ O	+ O ₃	3.00E-10	0.00	A42 FEHSENFELD ET AL (1969B)
108	O ₄ -	+ CO ₂	+ O ₂	1.00E-10	0.00	A43 ADAMS ET AL (1970)
109	O ₄ -	+ HV	+ CO ₄ -	4.30E-10	0.00	A44 PETERSON (1976)
110	CO ₃ -	+ HV	+ O ⁻	2.00E-01	0.00	A45 FERGUSON (1974)
111	CO ₃ -	+ O	+ O ₂	8.00E-11	0.00	A46 FEHSENFELD AND FERGUSON (1974)
112	CO ₃ -	+ NO ₂	+ NO ₃ -	2.00E-10	0.00	A47 BORTNER AND BAURER (1979)
113	CO ₄ -	+ HV	+ O ₂	1.40E-02	0.00	A48 WONG ET AL (1972)
114	CO ₄ -	+ HV	+ CO ₂	3.00E-01	0.00	A49 FEHSENFELD ET AL (1969B)
115	CO ₄ -	+ O	+ CO ₃ -	1.50E-10	0.00	A50 FEHSENFELD ET AL (1969B)
116	CO ₄ -	+ O ₃	+ CO ₂	1.30E-10	0.00	A51 FEHSENFELD ET AL (1969B)
117	CO ₄ -	+ NO	+ NO ₃ -	4.80E-11	0.00	A52 FEHSENFELD ET AL (1969B)
118	NO ₂ -	+ HV	+ NO ₂	3.00E-02	0.00	A53 BORTNER AND BAURER (1979)
119	NO ₂ -	+ O ₃	+ O ₂	1.80E-11	0.00	A54 FEHSENFELD AND FERGUSON (1968)
120	NO ₃ -	+ HV	+ O ₂	3.00E-02	0.00	A55 BORTNER AND BAURER (1979)
121	NO ₃ -	+ O	+ O ₂	1.00E-12	0.00	A56 FEHSENFELD ET AL (1969B)
122	NO ₃ -	+ NO	+ NO ₂	2.00E-11	0.00	A57 GRAHAM AND JOHNSTON (1978)
123	NO ₃ -	+ HV	+ NO ₃ -	2.00E-02	0.00	A58 BORTNER AND BAURER (1979)
124	O-	+ O ₃	+ O ₂	4.40E-10	0.00	A59 BORTNER AND BAURER (1979)
125	O ^{+(2D)}	+ N ₂	+ O	3.00E-10	0.00	A60 STEBBINGS ET AL (1966)
126	O ^{+(2D)}	+ O ₂	+ O ₂	3.00E-10	0.00	A61 STEBBINGS ET AL (1966)
127	O ⁺	+ CO ₂	+ CO	1.00E-09	0.00	A62 LINDBERGER ET AL (1974)
128	O	+ D	+ N	4.80E-33	-2.00	A63 CAMPBELL AND GRAY (1973)
129	O	+ O ₂	+ N	5.50E-34	-2.60	A64 KAUFMAN AND KELSO (1964)
130	O	+ O ₃	+ O ₂	1.20E-11	0.00	A65 WONG SCHIFF (1969)
131	O	+ O ₃	+ H	7.00E-11	0.00	A66 WONG AND POTTER (1965)
132	O	+ H ₂ 02	+ OH	1.20E-12	0.00	A67 DAVIS ET AL (1974A)
133	O	+ H ₂ 02	+ H ₂ O	1.80E-12	0.00	A68 DAVIS ET AL (1974A)
134	H	+ H ₂ 02	+ H ₂	2.30E-12	0.00	A69 KLEMm ET AL (1975)
135	H	+ H ₂ 02	+ OH	2.90E-12	0.00	A70 KLEMm ET AL (1975)
136	H	+ O ₂	+ M	4.40E-33	0.00	A71 BAULCH ET AL (1972)
137	N	+ O ₂	+ O	5.50E-12	0.00	A72 BECKER ET AL (1969)
138	N	+ NO	+ N ₂	8.20E-11	0.00	A73 CLYNE AND McDERMID (1975)
139	N	+ O	+ M	1.80E-31	-5.0	A74 BAULCH ET AL (1973)
140	OH	+ O ₃	+ O ₂	1.30E-12	0.00	A75 ANDERSON AND KAUFMAN (1973)
141	OH	+ H ₂ 02	+ H ₂ O	1.00E-02	-7.50E-02	A76 BORTNER AND BAURER (1979)

Table A1. Chemical Reaction Scheme (Contd)

NO.	CHEMICAL REACTION		REFERENCE
	A	B	
142	NO + O ₃ = O ₂	+ NO ₂	1.70E-12 0.00 -1.31E+03 452 BAULCH ET AL (1973)
143	NO + O = NO ₂	+ M	1.55E-32 0.00 -5.84E+02 452 MYTLOCK ET AL (1976)
144	H ₂ O ₂ + O = O ₂	+ H ₂ O ₂	7.55E-12 0.00 -5.00E+02 452 BORTNER AND BAURER (1979)
145	H ₂ O ₂ + O ₃ = O ₂	+ OH	1.00E-13 0.00 -1.25E+03 452 GARVIN (1973)
146	NO ₂ + H = OH	+ OH	4.80E-10 0.00 -4.00E+02 452 CLYNE AND MONKHOUSE (1977)
147	CO + OH = CO ₂	+ M	2.20E-13 0.00 -8.00E+01 452 DAVIS ET AL (1974B)
148	CO + O = CO ₂	+ M	6.50E-33 0.00 -2.18E+03 452 BAULCH ET AL (1976)
149	N ₂ O + E = N + HV	+ E	3.20E-10 0.00 0.00 BORTNER AND BAURER (1979)
150	NO ₂ + E = O + HV	+ O	2.00E-07 0.00 0.00 BORTNER AND BAURER (1979)
151	NO + E = O + HV	+ O	2.00E-07 0.00 0.00 BORTNER AND BAURER (1979)
152	O ₂ + E = O + O	+ O	2.00E-07 0.00 0.00 BORTNER AND BAURER (1979)
153	O ₂ + E = O + HV	+ O	3.50E-12 0.00 0.00 BORTNER AND BAURER (1979)
154	N ₂ + E = N + HV	+ N	3.50E-12 0.00 0.00 BORTNER AND BAURER (1979)
155	H ₂ + E = H + HV	+ H	3.50E-12 0.00 0.00 BORTNER AND BAURER (1979)
156	NO ₂ ⁺ + H ₂ O = NO + H ₂ O	+ H ₂ O	3.00E-07 0.00 0.00 BORTNER AND BAURER (1979)
157	NO ₂ ⁺ + H ₂ O = NO + H ₂ O	+ H ₂ O	1.50E-06 0.00 0.00 BORTNER AND BAURER (1979)
158	NO ₂ ⁺ + 2H ₂ O = NO + 3H ₂ O	+ H ₂ O	2.00E-06 0.00 0.00 BORTNER AND BAURER (1979)
159	NO ₂ ⁺ + 3H ₂ O = O ₂ + E	+ E	3.00E-06 0.00 0.00 BORTNER AND BAURER (1979)
160	O ₂ + E = O ₂ ⁺ + O	+ O ₂ ⁺	1.40E-29 0.00 -6.00E+02 452 PHELPS (1969)
161	O ₃ + E = O ₂ ⁺ + O ₂	+ O ₂ ⁺	9.00E-12 0.00 0.00 STELMAN ET AL (1972)
162	O ₂ ⁺ + O ₂ = O ₄ ⁺	+ O ₄ ⁺	3.90E-30 0.00 0.00 BORTNER AND BAURER (1979)
163	O ₄ ⁺ + O ₂ = O ₂ ⁺ + O ₂	+ O ₂ ⁺	2.00E-05 0.00 -5.40E+03 452 BORTNER AND BAURER (1979)
164	O ₂ ^{1D} + O ₂ = O ₂ ⁺ + O ₂	+ O ₂ ⁺	2.00E-05 0.00 -5.40E+03 452 BORTNER AND BAURER (1979)
165	NO ₂ + CO ₂ = N ₂ + NO ₃ ⁺	+ N ₂	3.00E-29 0.00 0.00 BORTNER AND BAURER (1979)
166	NO + H ₂ O = NO ₂ + H ₂ O	+ H ₂ O	1.50E-28 0.00 0.00 BORTNER AND BAURER (1979)
167	N ₂ ⁺ + O ₂ = O ₂ ⁺ + N ₂	+ N ₂	5.00E-11 0.00 0.00 BORTNER AND BAURER (1979)
168	O ₂ + N ₂ = NO ₂ + N	+ NO ₂	1.20E-12 0.00 0.00 LINDBERGER ET AL (1974)
169	O ₂ + O ₂ = O ₃ + O	+ O ₃	2.00E-11 0.00 0.00 LINDBERGER ET AL (1974)
170	H ₃ O ⁺ + H ₂ O = H ₅ O ₂ ⁺	+ H ₅ O ₂ ⁺	3.50E-27 0.00 0.00 GOOD ET AL (1970)
171	H ₃ O ⁺ + N ₂ = H ₃ O ⁺ .N ₂	+ N ₂	1.40E-30 0.00 0.00 BORTNER AND BAURER (1979)
172	H ₅ O ₂ ⁺ + H ₂ O = H ₇ O ₃ ⁺	+ H ₇ O ₃ ⁺	2.20E-27 0.00 0.00 BORTNER AND BAURER (1979)
173	H ₇ O ₃ ⁺ + H ₂ O = H ₉ O ₄ ⁺	+ H ₉ O ₄ ⁺	2.30E-27 0.00 0.00 BORTNER AND BAURER (1979)
174	H ₇ O ₃ ⁺ + H ₂ O = H ₅ O ₂ ⁺	+ H ₅ O ₂ ⁺	2.50E-03 0.00 -1.00E+04 452 BORTNER AND BAURER (1979)
175	H ₉ O ₄ ⁺ + H ₂ O = H ₇ O ₃ ⁺	+ H ₇ O ₃ ⁺	2.00E-01 0.00 -8.80E+03 452 BORTNER AND BAURER (1979)
176	H ₁₁ O ₅ ⁺ + H ₂ O = H ₉ O ₄ ⁺	+ H ₉ O ₄ ⁺	4.00E-01 0.00 -7.67E+03 452 BORTNER AND BAURER (1979)
177	NO ₂ ⁺ + H ₂ O = NO ₃ ⁺ + H ₂ O	+ NO ₃ ⁺	1.10E-27 0.00 0.00 BORTNER AND BAURER (1979)
178	NO ₂ ⁺ + 2H ₂ O = NO ₃ ⁺ + 3H ₂ O	+ NO ₃ ⁺	1.10E-27 0.00 0.00 BORTNER AND BAURER (1979)
179	O ₂ + O ₂ = O ₃ + O ₂	+ O ₃	1.10E-30 0.00 0.00 BORTNER AND BAURER (1979)
180	O ₂ + O ₂ = O ₄ ⁺	+ O ₄ ⁺	3.50E-31 0.00 0.00 BORTNER AND BAURER (1979)
181	O ₂ + O ₂ = CO ₂	+ CO ₂	2.00E-29 0.00 0.00 BORTNER AND BAURER (1979)
182	O ₃ + CO ₂ = CO ₃ ⁻	+ CO ₃ ⁻	5.50E-10 0.00 -4.49 0.00 BORTNER AND BAURER (1979)
183	O ₃ + N ₂ = NO ₂ + N ₂	+ NO ₂	2.00E-05 0.00 -6.30E+03 452 BORTNER AND BAURER (1979)
184	CO ₃ ⁻ + NO = NO ₂ + CO ₂	+ NO ₂	1.10E-11 0.00 0.00 BORTNER AND BAURER (1979)
185	O + HV = O ₂ + HV	+ O ₂	0.00 0.00 0.00 BORTNER AND BAURER (1979)
186	O ₂ + HV = O ₃ + HV	+ O ₃	0.00 0.00 0.00 BORTNER AND BAURER (1979)
187	N ₂ + HV = N ₂ ⁺ + HV	+ N ₂ ⁺	0.00 0.00 0.00 BORTNER AND BAURER (1979)
188	N ₂ ⁺ + HV = N ₂ + HV	+ N ₂	0.00 0.00 0.00 BORTNER AND BAURER (1979)

Wave length dependent (see text)

Table A1. Chemical Reaction Scheme (Contd)

NO.	CHEMICAL REACTION	REFERENCE		
		A	B	C
189 N	HV = N+	+ E	0.00	0.
190 H	HV = H+	+ E	0.00	0.
191 HE	HV = HE+	+ E	0.00	0.
192 O2	HV = O	+ + + O	0.00	0.
193 O2	HV = O1D	+ + O	0.00	0.
194 O3	HV = O2	+ + + O	0.00	0.
195 O3	HV = O21D	+ + O1D	0.00	0.
196 H2O	HV = OH	+ + H	0.00	0.
197 H2O2	HV = OH	+ + OH	0.00	0.
198 HO2	HV = O2	+ + H	0.00	0.
199 NO	HV = N	+ + O	0.00	0.
200 NO2	HV = NO	+ + O1D	0.00	0.
201 N2O	HV = N2	+ + O1D	0.00	0.
202 HNO2	HV = OH	+ + NO	0.00	0.
203 CO2	HV = CO	+ + O1D	0.00	0.
204 CO2	HV = CO	+ + O	0.00	0.
205 NO3	HV = NO + O2	+ + O2	0.00	0.
206 O	HV = O + (2D)	+ + E	0.00	0.
207 O21D	HV = O2+	+ + E	0.00	0.
208 E*	HV = N2	+ + N*	+ 2E	0.
209 E*	HV = O2	+ + O	+ 2E	0.
210 E*	HV = O2	+ + O	+ 2E	0.
211 E*	HV = O	+ + O	+ 2E	0.
212 E*	HV = O	+ + O	+ 2E	0.
213 E*	HV = N2	+ + N2+	+ 2E	0.
214 E*	HV = O2	+ + O2+	+ 2E	0.
215 E*	HV = N2	+ + N2D	+ E	0.00

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Appendix B

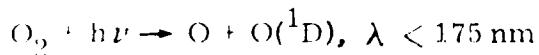
Photodissociation processes and Their Rate Coefficients

Appendix B contains information on the photodissociation, photoionization, and energetic electron impact reactions. Table B1 lists the solar fluxes in the wavelength region 116.2 to 730.0 nm along with the absorption cross sections for O₂, O₃, H₂O, H₂O₂, NO₂, NO₃, and N₂O. Table B2 contains the absorption cross sections for N₂O₅, HNO₂, HNO₃, CH₄, CO₂, CH₂O, and HO₂.

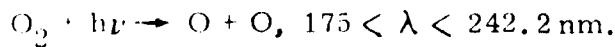
Table B3 lists the photodissociation processes along with their total photodissociation rate coefficient. The units of all the rate coefficients presented in this table and in all the tables that follow (Tables B4-B23) are sec⁻¹. The constants in Table B3 were computed from Eq. (52) using the entire unattenuated solar flux. The values listed here are indications of the maximum rate coefficient that each process can attain.

B1. MOLECULAR OXYGEN

Two distinct processes for O₂ photodissociation are considered,



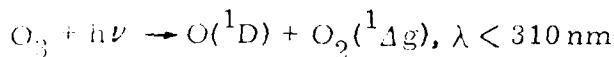
and



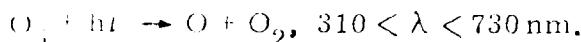
The cross sections for this molecule are taken from the compilation of Ackerman¹⁶ with the exception of those wavelengths in the Schumann Range bands. The photodissociation rate coefficients for the first process above are listed in Table B4, and those for the second process are listed in Table B5. The column densities given in these and all the other tables that follow (Tables B6-B23) are in units of cm^{-2} column.

B2. OZONE

For ozone, two processes are also considered,



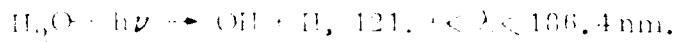
and



The absorption cross sections for ozone are also taken from the compilation of Ackerman.¹⁶ The photodissociation rate coefficients for the process with the excited products are listed in Table B6, and those for the process with the ground state products are listed in Table B7.

B3. WATER VAPOR

Water vapor is photodissociated primarily through the process

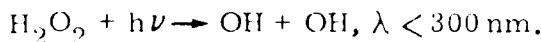


The absorption cross sections for this molecule are taken from Watanabe and Zelikoff.¹⁷ The photodissociation rate coefficients for this process are tabulated in Table B7.

¹⁶ Ackerman, R., and Zelikoff, M.J. (1953) Absorption coefficients of water vapor in the vacuum ultraviolet, J. Opt. Soc. Am. 43:753-755.

B1. HYDROGEN PEROXIDE

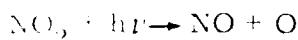
According to Urey et al.^{B2} and Holt et al.,^{B3} the photodissociation process for hydrogen peroxide is probably



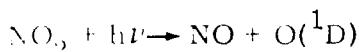
The absorption cross sections above 225 nm are from Urey et al., and those below 225 nm are from Holt et al. The photodissociation rate coefficients for this molecule are presented in Table B9.

B5. NITROGEN DIOXIDE

The primary quantum yield for nitrogen dioxide is unity for the process



for wavelengths from 398 to 244 nm. The process



has unity primary quantum yield from 244 to 190 nm. For these computations, however, all of the photodissociation of NO_2 is assumed to take place according to the latter process. The absorption cross sections for this molecule are taken from Nakayama et al.^{B4} and from Hall and Blacet.^{B5} The photodissociation rate coefficients for this process are listed in Table B10.

B6. THE NO_3 RADICAL

According to Johnston and Graham,¹³ the photolysis of NO_3 can proceed as

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- B2. Urey, H.C., Dawson, L.H., and Rice, F.O. (1929) The absorption spectrum and decomposition of hydrogen peroxide by light, *J. Am. Chem. Soc.*, 51:1371-1383.
 - B3. Holt, R.B., McElveen, C.K., and Oldenberg, O. (1943) Ultraviolet absorption spectrum of hydrogen peroxide, *J. Chem. Phys.*, 16:225-229.
 - B4. Nakayama, T., Kitamura, M.V., and Watanabe, K. (1959) Ionization potential and absorption coefficients of nitrogen dioxide, *J. Chem. Phys.*, 30:1115-1126.
 - B5. Hall, T.C., and Blacet, E.P. (1952) Separation of the absorption spectra of NO_2 and N_2O_4 in the region 2400-3000 Å, *J. Chem. Phys.*, 20:1745-1751.

Table B2. Total photoionization Rate Coefficients for the Process $O_2 + h\nu \rightarrow O + O(^1D)$ for the Wavelengths Below 1 μ m (Cont'd)

	03 COLUMN DENSITY									
	2.0E+15	5.0E+15	1.0E+16	2.0E+16	5.0E+16	1.0E+17	2.0E+17	5.0E+17	1.0E+18	2.0E+18
1.0E+13	1.66806E-06	1.65171E-06	1.62513E-06	1.57431E-06	1.43824E-06	1.25387E-06	9.88396E-07	5.71980E-07	2.94155E-07	1.02772E-07
2.0E+13	1.66793E-06	1.65158E-06	1.62500E-06	1.57419E-06	1.43814E-06	1.25378E-06	9.88334E-07	5.71953E-07	2.94145E-07	1.02770E-07
5.0E+13	1.66754E-06	1.6519E-06	1.62462E-06	1.57383E-06	1.43782E-06	1.25351E-06	9.88149E-07	5.71873E-07	2.94115E-07	1.02762E-07
1.0E+14	1.66688E-06	1.65055E-06	1.62399E-06	1.57322E-06	1.43728E-06	1.25307E-06	9.87839E-07	5.71740E-07	2.94066E-07	1.02750E-07
2.0E+14	1.66558E-06	1.64926E-06	1.62273E-06	1.57201E-06	1.43621E-06	1.25219E-06	9.87221E-07	5.71474E-07	2.93967E-07	1.02725E-07
5.0E+14	1.66167E-06	1.64500E-06	1.61895E-06	1.56839E-06	1.43302E-06	1.24956E-06	9.85369E-07	5.70676E-07	2.93670E-07	1.02651E-07
1.0E+15	1.65518E-06	1.63899E-06	1.61268E-06	1.56238E-06	1.42771E-06	1.24519E-06	9.82294E-07	5.69351E-07	2.93177E-07	1.02528E-07
2.0E+15	1.64230E-06	1.62629E-06	1.60324E-06	1.55047E-06	1.41718E-06	1.23652E-06	9.76187E-07	5.66715E-07	2.92195E-07	1.02282E-07
5.0E+15	1.60450E-06	1.58897E-06	1.56372E-06	1.51546E-06	1.38623E-06	1.21102E-06	9.58212E-07	5.58929E-07	2.89282E-07	1.01550E-07
1.0E+16	1.54414E-06	1.52939E-06	1.50540E-06	1.45956E-06	1.33677E-06	1.17020E-06	9.29361E-07	5.46341E-07	2.84538E-07	1.00349E-07
2.0E+16	1.43268E-06	1.41934E-06	1.39766E-06	1.35621E-06	1.24522E-06	1.09445E-06	8.75550E-07	5.22546E-07	2.75441E-07	9.80170E-08
5.0E+16	1.1604CE-06	1.15043E-06	1.13421E-06	1.10323E-06	1.02023E-06	9.07081E-07	7.40652E-07	4.60789E-07	2.50972E-07	9.15433E-08
1.0E+17	8.54413E-07	8.47991E-07	8.37554E-07	8.17620E-07	7.64191E-07	6.90938E-07	5.80801E-07	3.82530E-07	2.17829E-07	8.22435E-08
2.0E+17	5.36644E-07	5.33460E-07	5.28291E-07	5.18431E-07	4.92011E-07	4.55440E-07	3.98265E-07	2.82798E-07	1.70978E-07	6.78314E-08
5.0E+17	2.54983E-07	2.53852E-07	2.52026E-07	2.48539E-07	2.39264E-07	2.26352E-07	2.05261E-07	1.56698E-07	1.01221E-07	4.28533E-08
1.0E+18	1.36289E-07	1.35699E-07	1.34746E-07	1.32943E-07	1.28121E-07	1.21616E-07	1.10911E-07	8.58965E-08	5.64943E-08	2.44971E-08
2.0E+18	6.00193E-08	5.37393E-08	5.92891E-08	5.84436E-08	5.62472E-08	5.32901E-08	4.85643E-08	3.76423E-08	2.48191E-08	1.08074E-08
5.0E+18	1.13865E-08	1.13204E-08	1.12151E-08	1.10204E-08	1.05331E-08	9.91549E-09	8.98713E-09	6.92906E-09	4.54882E-09	1.96587E-09
1.0E+19	1.66447E-09	1.65246E-09	1.63346E-09	1.59880E-09	1.51493E-09	1.41452E-09	1.27268E-09	9.73832E-10	6.36991E-10	2.74164E-10
2.0E+19	9.58961E-11	9.48877E-11	9.32951E-11	9.04033E-11	8.35107E-11	7.55905E-11	6.54293E-11	4.79128E-11	3.1048E-11	1.33560E-11
5.0E+19	1.95990E-12	1.91293E-12	1.83731E-12	1.69528E-12	1.33423E-12	9.00690E-13	4.20783E-13	5.68626E-14	9.60759E-15	3.42372E-15
1.0E+20	4.57705E-13	4.46212E-13	4.27722E-13	3.93016E-13	3.05030E-13	2.00223E-13	8.67228E-14	7.32785E-15	1.33288E-16	7.01575E-20
2.0E+20	4.61935E-14	4.50798E-14	4.32836E-14	3.99043E-14	3.2770E-14	2.08583E-14	9.30490E-15	6.43112E-16	1.61428E-17	6.38520E-21
5.0E+20	9.31927E-17	9.10317E-17	8.75410E-17	8.09560E-17	6.40283E-17	4.33116E-17	1.98226E-17	1.57073E-20	1.84043E-18	3.84043E-23
1.0E+21	4.18259E-21	4.08390E-21	3.92769E-21	3.63297E-21	2.87448E-21	1.94652E-21	8.92285E-22	8.59497E-23	1.73975E-24	7.12828E-28
2.0E+21	8.61614E-30	8.41687E-30	8.09493E-30	7.48752E-30	5.92534E-30	4.01179E-30	1.83903E-30	1.77149E-31	3.58584E-33	1.46924E-36
5.0E+21	7.54474E-56	7.37024E-56	7.08833E-56	6.51646E-56	5.18853E-56	3.51293E-56	1.61035E-56	1.55121E-57	3.13995E-59	1.286555E-62
1.3E+22	2.80670E-99	2.74178E-99	2.63691E-99	2.43905E-99	2.193017E-99	1.30684E-99	5.99062E-100	5.77062E-101	1.16808E-102	4.78605E-105

Table 11. Total Photoionization Rate Coefficients for the Process $O_2 + h\nu \rightarrow O + O(^1D)$ for the Wavelengths Below 175 nm

	O3 COLUMN DENSITY						
	1.0E+12	2.0E+12	5.0E+12	1.0E+13	2.0E+13	5.0E+13	1.0E+14
1.0E+13	1.67913E-06	1.67912E-06	1.67908E-06	1.67902E-06	1.67885E-06	1.67858E-06	1.67358E-06
2.0E+13	1.67905E-06	1.67899E-06	1.67895E-06	1.67889E-06	1.67844E-06	1.67872E-06	1.67345E-06
5.0E+13	1.67860E-06	1.67857E-06	1.67855E-06	1.67849E-06	1.67833E-06	1.67805E-06	1.67305E-06
1.0E+14	1.67794E-06	1.67793E-06	1.6777E-06	1.67783E-06	1.67767E-06	1.67739E-06	1.67517E-06
2.0E+14	1.67662E-06	1.67662E-06	1.67660E-06	1.67657E-06	1.67635E-06	1.67607E-06	1.67109E-06
5.0E+14	1.67268E-06	1.67267E-06	1.67263E-06	1.67257E-06	1.67241E-06	1.67213E-06	1.66716E-06
1.0E+15	1.66613E-06	1.66613E-06	1.66611E-06	1.66608E-06	1.66603E-06	1.66586E-06	1.66533E-06
2.0E+15	1.65314E-06	1.65314E-06	1.65312E-06	1.65309E-06	1.65304E-06	1.65286E-06	1.65042E-06
5.0E+15	1.61531E-06	1.61531E-06	1.61501E-06	1.61499E-06	1.61496E-06	1.61475E-06	1.64771E-06
1.0E+16	1.55413E-06	1.55413E-06	1.55411E-06	1.55409E-06	1.55404E-06	1.55388E-06	1.55163E-06
2.0E+16	1.44170E-06	1.44170E-06	1.44168E-06	1.44166E-06	1.44148E-06	1.44125E-06	1.43944E-06
5.0E+16	1.16714E-06	1.16715E-06	1.16714E-06	1.16712E-06	1.16709E-06	1.16699E-06	1.16377E-06
1.0E+17	8.58763E-07	8.58761E-07	8.58754E-07	8.58743E-07	8.58721E-07	8.58656E-07	8.58546E-07
2.0E+17	5.38802E-07	5.38800E-07	5.38797E-07	5.38792E-07	5.38781E-07	5.38748E-07	5.38694E-07
5.0E+17	2.55751E-07	2.55751E-07	2.55750E-07	2.55749E-07	2.55744E-07	2.55732E-07	2.55713E-07
1.0E+18	1.35653E-07	1.35650E-07	1.35649E-07	1.35649E-07	1.36686E-07	1.36680E-07	1.36489E-07
2.0E+18	6.02101E-08	6.02098E-08	6.02098E-08	6.02098E-08	6.02055E-08	6.02006E-08	6.01910E-08
5.0E+18	1.14317E-08	1.14317E-08	1.14316E-08	1.14315E-08	1.14306E-08	1.14272E-08	1.14203E-08
1.0E+19	1.67274E-09	1.67273E-09	1.67272E-09	1.67270E-09	1.67266E-09	1.67190E-09	1.67165E-09
2.0E+19	9.65914E-11	9.65911E-11	9.65900E-11	9.65883E-11	9.65742E-11	9.65565E-11	9.62415E-11
5.0E+19	1.99185E-12	1.99183E-12	1.99178E-12	1.99170E-12	1.99154E-12	1.99106E-12	1.98864E-12
1.0E+20	4.65523E-13	4.65519E-13	4.65508E-13	4.65488E-13	4.65448E-13	4.65330E-13	4.63559E-13
2.0E+20	4.69508E-14	4.69504E-14	4.69493E-14	4.69474E-14	4.69436E-14	4.69321E-14	4.67606E-14
5.0E+20	9.46611E-17	9.46603E-17	9.46581E-17	9.46544E-17	9.46470E-17	9.46248E-17	9.45139E-17
1.0E+21	4.24629E-21	4.24626E-21	4.24616E-21	4.24594E-21	4.24467E-21	4.23971E-21	4.22980E-21
2.0E+21	8.75154E-30	8.75147E-30	8.75127E-30	8.75093E-30	8.74479E-30	8.73797E-30	8.68361E-30
5.0E+21	7.66330E-56	7.66324E-56	7.66306E-56	7.66276E-56	7.66216E-56	7.63353E-56	7.63141E-56
1.0E+22	2.85080E-99	2.85078E-99	2.85060E-99	2.85038E-99	2.84972E-99	2.84860E-99	2.82868E-99

Table B3. Total Photodissociation Rate Coefficients for Unattenuated Solar Flux

$O_2 + HV = O + O_1D$	$1.67927E-06 \text{ SEC}^{-1}$
$O_2 + HV = O + O$	$8.36798E-08 \text{ SEC}^{-1}$
$O_3 + HV = O_1D + O_2ID$	$9.72665E-03 \text{ SEC}^{-1}$
$O_3 + HV = O + O_2$	$4.17996E-04 \text{ SEC}^{-1}$
$H_2O + HV = OH + H$	$7.38662E-06 \text{ SEC}^{-1}$
$H_2O_2 + HV = OH + OH$	$1.59387E-04 \text{ SEC}^{-1}$
$NO_2 + HV = NO + O_1D$	$1.51039E-02 \text{ SEC}^{-1}$
$NO_3 + HV = NO + O_2$	$7.04137E-02 \text{ SEC}^{-1}$
$N_2O + HV = N_2 + O_1D$	$7.76688E-07 \text{ SEC}^{-1}$
$N_2O_5 + HV = NO_2 + NO_2$	$6.92337E-04 \text{ SEC}^{-1}$
$HNO_2 + HV = OH + NO$	$5.84756E-04 \text{ SEC}^{-1}$
$HNO_3 + HV = OH + NO_2$	$1.16540E-04 \text{ SEC}^{-1}$
$CH_4 + HV = CH_3 + H$	$9.08606E-06 \text{ SEC}^{-1}$
$CO_2 + HV = CO + O_1D$	$8.84043E-08 \text{ SEC}^{-1}$
$CO_2 + HV = CO + O$	$8.24894E-09 \text{ SEC}^{-1}$
$CH_2O + HV = CHO + H$	$2.58902E-04 \text{ SEC}^{-1}$
$HO_2 + HV = O_2 + H$	$1.13657E-03 \text{ SEC}^{-1}$

Table B2. Absorption Cross Sections for N₂O₅, HNO₂, HNO₃, CH₄, CO₂, CH₂O, and HO₂ (Contd)

WAVELENGTH BAND ANGSTROMS	N205	HNO2	CROSS SECTIONS (CM ²)				
			HNO3	CH4	CO2	CH2O	
3200.00-3250.00	6.00E-21	2.00E-20	1.00E-22	0.	0.	2.20E-20	0.
3250.00-3300.00	5.00E-21	2.00E-20	0.	0.	0.	3.40E-20	0.
3300.00-3350.00	3.40E-21	3.00E-20	0.	0.	0.	1.50E-20	0.
3350.00-3400.00	3.00E-21	2.80E-20	0.	0.	0.	1.20E-20	0.
3400.00-3450.00	2.40E-21	3.10E-20	0.	0.	0.	2.00E-20	0.
3450.00-3500.00	2.00E-21	3.00E-20	0.	0.	0.	5.00E-21	0.
3500.00-3550.00	1.50E-21	5.00E-20	0.	0.	0.	2.80E-21	0.
3550.00-3600.00	1.20E-21	5.00E-20	0.	0.	0.	1.20E-21	0.
3600.00-3650.00	9.00E-22	2.90E-20	0.	0.	0.	0.	0.
3650.00-3700.00	7.00E-22	8.00E-20	0.	0.	0.	0.	0.
3700.00-3750.00	4.00E-22	4.00E-20	0.	0.	0.	0.	0.
3750.00-3800.00	2.50E-22	2.00E-20	0.	0.	0.	0.	0.
3800.00-3850.00	0.	4.00E-20	0.	0.	0.	0.	0.
3850.00-3900.00	0.	5.00E-20	0.	0.	0.	0.	0.
3900.00-3950.00	0.	0.	0.	0.	0.	0.	0.
3950.00-4000.00	0.	0.	0.	0.	0.	0.	0.
4000.00-4050.00	0.	0.	0.	0.	0.	0.	0.
4050.00-4100.00	0.	0.	0.	0.	0.	0.	0.
4100.00-4150.00	0.	0.	0.	0.	0.	0.	0.
4150.00-4200.00	0.	0.	0.	0.	0.	0.	0.
4200.00-4250.00	0.	0.	0.	0.	0.	0.	0.
4250.00-4300.00	0.	0.	0.	0.	0.	0.	0.
4300.00-4350.00	0.	0.	0.	0.	0.	0.	0.
4350.00-4400.00	0.	0.	0.	0.	0.	0.	0.
4400.00-4450.00	0.	0.	0.	0.	0.	0.	0.
4450.00-4500.00	0.	0.	0.	0.	0.	0.	0.
4500.00-4550.00	0.	0.	0.	0.	0.	0.	0.
4550.00-4600.00	0.	0.	0.	0.	0.	0.	0.
4600.00-4650.00	0.	0.	0.	0.	0.	0.	0.
4650.00-4700.00	0.	0.	0.	0.	0.	0.	0.
4700.00-4750.00	0.	0.	0.	0.	0.	0.	0.
4750.00-4800.00	0.	0.	0.	0.	0.	0.	0.
4800.00-4850.00	0.	0.	0.	0.	0.	0.	0.
4850.00-4900.00	0.	0.	0.	0.	0.	0.	0.
4900.00-4950.00	0.	0.	0.	0.	0.	0.	0.
4950.00-5000.00	0.	0.	0.	0.	0.	0.	0.
5000.00-5050.00	0.	0.	0.	0.	0.	0.	0.
5050.00-5100.00	0.	0.	0.	0.	0.	0.	0.
5100.00-5150.00	0.	0.	0.	0.	0.	0.	0.
5150.00-5200.00	0.	0.	0.	0.	0.	0.	0.
5200.00-5250.00	0.	0.	0.	0.	0.	0.	0.
5250.00-5300.00	0.	0.	0.	0.	0.	0.	0.
5300.00-5350.00	0.	0.	0.	0.	0.	0.	0.
5350.00-5400.00	0.	0.	0.	0.	0.	0.	0.
5400.00-5450.00	0.	0.	0.	0.	0.	0.	0.
5450.00-5500.00	0.	0.	0.	0.	0.	0.	0.
5500.00-5550.00	0.	0.	0.	0.	0.	0.	0.
5550.00-5600.00	0.	0.	0.	0.	0.	0.	0.
5600.00-5650.00	0.	0.	0.	0.	0.	0.	0.
5650.00-5700.00	0.	0.	0.	0.	0.	0.	0.
5700.00-5750.00	0.	0.	0.	0.	0.	0.	0.
5750.00-5800.00	0.	0.	0.	0.	0.	0.	0.

Table B2. Absorption Cross Sections for N₂O₅, HNO₂, HNO₃, CH₄, CO₂, CH₂O, and HO₂ (Contd)

WAVELENGTH BAND ANGSTROMS	N2O5	HNO2	HNO3	CH4	C02	CH2O	HO2
1792.60-1803.60	0.	0.	0.	1.30E-21	0.	5.50E-18	
1803.60-1816.40	0.	0.	0.	1.00E-21	0.	5.50E-18	
1816.40-1830.60	0.	0.	0.	6.00E-22	0.	5.50E-18	
1830.60-1846.20	0.	0.	0.	4.00E-22	0.	5.50E-18	
1846.20-1863.40	0.	0.	0.	3.00E-22	0.	5.60E-18	
1863.40-1882.20	0.	0.	0.	1.50E-22	0.	5.70E-18	
1882.20-1902.40	0.	0.	0.	8.00E-23	0.	5.80E-18	
1902.40-1924.00	0.	0.	0.	3.00E-23	0.	5.80E-18	
1924.00-1947.00	0.	0.	0.	2.00E-23	0.	5.80E-18	
1947.00-1971.80	0.	0.	0.	1.00E-23	0.	5.80E-18	
1971.80-1985.00	0.	8.00E-18	0.	8.00E-24	0.	5.80E-18	
1985.00-2000.00	0.	7.00E-18	0.	5.50E-24	0.	5.90E-18	
2000.00-2025.00	0.	5.00E-18	0.	4.00E-24	0.	5.90E-18	
2025.00-2050.00	0.	3.80E-18	0.	3.20E-24	0.	5.90E-18	
2050.00-2061.86	0.	3.00E-18	0.	3.00E-24	0.	5.90E-18	
2061.86-2083.33	0.	2.40E-18	0.	3.00E-24	0.	6.00E-18	
2083.33-2105.26	0.	1.70E-18	0.	2.80E-24	0.	6.00E-18	
2105.26-2127.66	5.00E-18	0.	6.00E-19	0.	2.40E-24	0.	6.00E-18
2127.68-2150.54	4.00E-18	0.	4.60E-19	0.	2.00E-24	0.	6.00E-18
2150.54-2173.91	3.30E-18	0.	3.40E-19	0.	0.	0.	5.80E-18
2173.91-2197.80	2.80E-18	0.	2.40E-19	0.	0.	0.	5.50E-18
2197.80-2222.22	2.20E-18	0.	1.50E-19	0.	0.	0.	5.00E-18
2222.22-2247.19	2.00E-18	0.	1.10E-19	0.	0.	0.	4.50E-18
2247.19-2272.73	1.50E-18	0.	9.00E-20	0.	0.	0.	4.00E-18
2272.73-2298.85	1.10E-18	0.	8.00E-20	0.	0.	0.	3.70E-18
2298.85-2325.58	1.00E-18	0.	5.50E-20	0.	0.	0.	3.00E-18
2325.58-2352.94	9.00E-19	0.	4.50E-20	0.	0.	0.	2.70E-18
2352.94-2380.95	8.00E-19	0.	3.80E-20	0.	0.	0.	2.30E-18
2380.95-2409.64	7.00E-19	0.	3.00E-20	0.	0.	0.	2.10E-18
2409.64-2439.02	6.00E-19	0.	2.80E-20	0.	0.	3.50E-21	1.70E-18
2439.02-2469.14	5.00E-19	0.	2.10E-20	0.	0.	3.80E-21	1.20E-18
2469.14-2500.00	4.00E-19	0.	2.00E-20	0.	0.	4.00E-21	1.00E-18
2500.00-2531.85	3.80E-19	0.	2.00E-20	0.	0.	6.00E-21	5.00E-19
2531.65-2564.10	3.20E-19	0.	2.00E-20	0.	0.	7.00E-21	3.80E-19
2564.10-2597.40	3.00E-19	0.	2.00E-20	0.	0.	8.00E-21	2.60E-19
2597.40-2631.58	2.40E-19	0.	2.00E-20	0.	0.	9.00E-21	1.10E-19
2631.58-2666.67	2.10E-19	0.	2.00E-20	0.	0.	1.00E-20	8.00E-20
2666.67-2702.70	1.80E-19	0.	2.00E-20	0.	0.	1.20E-20	5.00E-20
2702.70-2739.73	1.50E-19	0.	1.80E-20	0.	0.	1.50E-20	2.20E-20
2739.73-2777.78	1.30E-19	0.	1.60E-20	0.	0.	1.70E-20	0.
2777.78-2816.90	1.00E-19	0.	1.20E-20	0.	0.	1.90E-20	0.
2816.90-2857.14	8.00E-20	0.	1.10E-20	0.	0.	2.10E-20	0.
2857.14-2898.55	6.00E-20	0.	9.00E-21	0.	0.	2.50E-20	0.
2898.55-2941.18	4.00E-20	0.	6.50E-21	0.	0.	3.00E-20	0.
2941.18-2985.07	3.30E-20	0.	5.00E-21	0.	0.	3.80E-20	0.
2985.07-3030.30	2.80E-20	7.00E-21	3.20E-21	0.	0.	3.00E-20	0.
3030.30-3076.92	2.10E-20	7.50E-21	2.30E-21	0.	0.	4.00E-20	0.
3076.92-3100.00	1.90E-20	9.00E-21	1.80E-21	0.	0.	3.50E-20	0.
3100.00-3150.00	1.10E-20	1.00E-20	7.00E-22	0.	0.	4.50E-20	0.
3150.00-3200.00	9.00E-21	1.10E-20	3.90E-22	0.	0.	3.00E-20	0.

Table B2. Absorption Cross Sections for N_2O_5 , HNO_2 , HNO_3 , CH_4 , CO_2 , CH_2O , and HO_2

WAVELENGTH BAND ANGSTROMS	CROSS SECTIONS (CM ²)						
	N_2O_5	HNO_2	HNO_3	CH_4	CO_2	CH_2O	HO_2
1215.67-1215.67 0.	0.	0.	2.80E-17	7.30E-20	9.00E-18	0.	0.
1162.79-1169.59 0.	0.	0.	2.80E-17	1.10E-19	0.	0.	0.
1169.59-1176.47 0.	0.	0.	2.80E-17	8.80E-20	0.	0.	0.
1176.47-1183.43 0.	0.	0.	2.80E-17	6.00E-20	0.	0.	0.
1183.43-1190.48 0.	0.	0.	2.80E-17	4.70E-20	0.	0.	0.
1190.48-1197.60 0.	0.	0.	2.80E-17	3.80E-20	0.	0.	0.
1197.60-1204.82 0.	0.	0.	2.80E-17	4.00E-20	0.	0.	0.
1204.82-1212.12 0.	0.	0.	2.80E-17	6.10E-20	0.	0.	0.
1212.12-1219.51 0.	0.	0.	2.80E-17	7.40E-20	0.	0.	0.
1219.51-1226.99 0.	0.	0.	2.80E-17	1.00E-19	0.	0.	0.
1226.99-1234.57 0.	0.	0.	2.80E-17	1.20E-19	0.	0.	0.
1234.57-1242.24 0.	0.	0.	2.80E-17	1.60E-19	0.	0.	0.
1242.24-1250.00 0.	0.	0.	2.80E-17	1.80E-19	0.	0.	0.
1250.00-1257.86 0.	0.	0.	2.80E-17	2.10E-19	0.	0.	0.
1257.86-1265.82 0.	0.	0.	2.80E-17	2.80E-19	0.	0.	0.
1265.82-1273.89 0.	0.	0.	2.80E-17	3.30E-19	0.	0.	0.
1273.89-1282.05 0.	0.	0.	2.80E-17	4.20E-19	0.	0.	0.
1282.05-1290.32 0.	0.	0.	2.80E-17	5.90E-19	0.	0.	0.
1290.32-1298.70 0.	0.	0.	2.80E-17	7.00E-19	0.	0.	0.
1298.70-1307.19 0.	0.	0.	2.47E-17	8.00E-19	0.	0.	0.
1307.19-1315.79 0.	0.	0.	2.24E-17	8.30E-19	0.	0.	0.
1315.79-1324.50 0.	0.	0.	2.07E-17	8.50E-19	0.	0.	0.
1324.50-1333.33 0.	0.	0.	1.57E-17	8.70E-19	0.	0.	0.
1333.33-1342.28 0.	0.	0.	1.33E-17	8.70E-19	0.	0.	0.
1342.28-1351.35 0.	0.	0.	1.20E-17	8.70E-19	0.	0.	0.
1351.35-1360.54 0.	0.	0.	1.00E-17	8.60E-19	0.	0.	0.
1360.54-1369.86 0.	0.	0.	8.40E-18	8.10E-19	0.	0.	0.
1369.86-1379.31 0.	0.	0.	5.90E-18	7.20E-19	0.	0.	0.
1379.31-1388.89 0.	0.	0.	3.93E-18	6.40E-19	0.	0.	0.
1388.89-1408.45 0.	0.	0.	2.24E-18	6.00E-19	0.	0.	0.
1408.45-1428.57 0.	0.	0.	6.72E-19	5.70E-19	1.00E-17	0.	0.
1428.57-1449.28 0.	0.	0.	6.72E-20	5.50E-19	7.00E-18	0.	0.
1449.28-1470.59 0.	0.	0.	1.68E-20	5.70E-19	4.00E-18	0.	0.
1470.59-1492.54 0.	0.	0.	1.12E-20	5.70E-19	5.50E-18	0.	0.
1492.54-1515.15 0.	0.	0.	9.00E-21	5.00E-19	1.20E-17	0.	0.
1515.15-1538.46 0.	0.	0.	5.60E-21	4.00E-19	1.80E-17	0.	0.
1538.46-1562.50 0.	0.	0.	4.48E-21	3.00E-19	9.00E-18	0.	0.
1562.50-1587.30 0.	0.	0.	2.80E-21	2.50E-19	8.00E-19	0.	0.
1587.30-1612.90 0.	0.	0.	1.12E-21	1.60E-19	0.	0.	0.
1612.90-1639.34 0.	0.	0.	8.00E-22	1.00E-19	0.	0.	0.
1639.34-1666.67 0.	0.	0.	0.	7.00E-20	0.	0.	0.
1666.67-1694.92 0.	0.	0.	0.	4.00E-20	1.50E-18	0.	0.
1694.92-1724.14 0.	0.	0.	0.	1.50E-20	1.20E-17	0.	0.
1724.14-1739.13 0.	0.	0.	0.	1.00E-20	6.00E-18	0.	0.
1739.13-1750.00 0.	0.	0.	0.	8.00E-21	1.80E-18	0.	0.
1750.00-1763.20 0.	0.	0.	0.	5.00E-21	0.	0.	0.
1763.20-1768.60 0.	0.	0.	0.	3.50E-21	0.	0.	0.
1768.60-1774.80 0.	0.	0.	0.	3.00E-21	0.	0.	0.
1774.80-1781.60 0.	0.	0.	0.	2.00E-21	0.	0.	0.
1781.60-1792.60 0.	0.	0.	0.	1.50E-21	0.	0.	0.

Table B1. Solar Fluxes and Absorption Cross Sections for O₂, O₃, H₂O, H₂O₂, NO₂, NO₃, and N₂O (Contd)

WAVELENGTH BAND ANGSTROMS	FLUX PH/CM ² /S/A	CROSS SECTIONS (CM ²)						
		O2	O3	H2O	H2O2	NO2	NO3	N2O
5700.00-5750.00	5.38E+13 0.		4.75E-21 0.	0.	0.	6.00E-19 0.		
5750.00-5800.00	5.42E+13 0.		4.55E-21 0.	0.	0.	7.00E-19 0.		
5800.00-5850.00	5.42E+13 0.		4.35E-21 0.	0.	0.	7.00E-19 0.		
5850.00-5900.00	5.42E+13 0.		4.42E-21 0.	0.	0.	9.00E-19 0.		
5900.00-5950.00	5.44E+13 0.		4.61E-21 0.	0.	0.	1.00E-18 0.		
5950.00-6000.00	5.44E+13 0.		4.89E-21 0.	0.	0.	9.00E-19 0.		
6000.00-6050.00	5.42E+13 0.		4.84E-21 0.	0.	0.	8.00E-19 0.		
6050.00-6100.00	5.40E+13 0.		4.54E-21 0.	0.	0.	7.00E-19 0.		
6100.00-6150.00	5.40E+13 0.		4.24E-21 0.	0.	0.	5.00E-19 0.		
6150.00-6200.00	5.40E+13 0.		3.90E-21 0.	0.	0.	4.00E-19 0.		
6200.00-6250.00	5.38E+13 0.		3.60E-21 0.	0.	0.	2.00E-18 0.		
6250.00-6300.00	5.36E+13 0.		3.43E-21 0.	0.	0.	1.80E-18 0.		
6300.00-6350.00	5.34E+13 0.		3.17E-21 0.	0.	0.	6.00E-19 0.		
6350.00-6400.00	5.32E+13 0.		2.74E-21 0.	0.	0.	4.00E-19 0.		
6400.00-6450.00	5.30E+13 0.		2.61E-21 0.	0.	0.	2.20E-19 0.		
6450.00-6500.00	7.90E+13 0.		2.40E-21 0.	0.	0.	2.00E-19 0.		
6500.00-6600.00	5.22E+13 0.		2.07E-21 0.	0.	0.	3.00E-19 0.		
6600.00-6700.00	5.18E+13 0.		1.72E-21 0.	0.	0.	2.00E-18 0.		
6700.00-6800.00	5.14E+13 0.		1.37E-21 0.	0.	0.	2.00E-19 0.		
6800.00-6900.00	5.09E+13 0.		1.11E-21 0.	0.	0.	0.	0.	
6900.00-7000.00	5.04E+13 0.		9.13E-22 0.	0.	0.	0.	0.	
7000.00-7100.00	4.99E+13 0.		7.93E-22 0.	0.	0.	0.	0.	
7100.00-7200.00	4.94E+13 0.		6.40E-22 0.	0.	0.	0.	0.	
7200.00-7300.00	4.90E+13 0.		5.14E-22 0.	0.	0.	0.	0.	

Table B1. Solar Fluxes and Absorption Cross Sections for O₂, O₃, H₂O, H₂O₂, NO₂, NO₃, and N₂O (Contd)

WAVELENGTH BAND ANGSTROMS	FLUX PH/CM ² /S, A	CROSS SECTIONS (CM ²)						
		O2	O3	H2O	H2O2	NO2	NO3	N2O
3200.00-3250.00	1.62E+13 0.		1.50E-20 0.		0.	2.94E-19 0.		0.
3250.00-3300.00	1.94E+13 0.		7.78E-21 0.		0.	3.38E-19 0.		0.
3300.00-3350.00	1.79E+13 0.		3.72E-21 0.		0.	3.61E-19 0.		0.
3350.00-3400.00	1.89E+13 0.		1.71E-21 0.		0.	4.17E-19 0.		0.
3400.00-3450.00	2.02E+13 0.		7.46E-22 0.		0.	4.30E-19 0.		0.
3450.00-3500.00	2.06E+13 0.		2.66E-22 0.		0.	4.48E-19 0.		0.
3500.00-3550.00	2.06E+13 0.		1.09E-22 0.		0.	5.53E-19 0.		0.
3550.00-3600.00	2.08E+13 0.		5.49E-23 0.		0.	4.96E-19 0.		0.
3600.00-3650.00	2.36E+13 0.		0.		0.	5.84E-19 0.		0.
3650.00-3700.00	2.46E+13 0.		0.		0.	5.70E-19 0.		0.
3700.00-3750.00	2.48E+13 0.		0.		0.	5.74E-19 0.		0.
3750.00-3800.00	2.34E+13 0.		0.		0.	6.52E-19 0.		0.
3800.00-3850.00	2.22E+13 0.		0.		0.	6.45E-19 0.		0.
3850.00-3900.00	2.18E+13 0.		0.		0.	6.51E-19 0.		0.
3900.00-3950.00	2.38E+13 0.		0.		0.	6.08E-19 0.		0.
3950.00-4000.00	3.08E+13 0.		0.		0.	6.68E-19 0.		0.
4000.00-4050.00	3.80E+13 0.		0.		0.	6.52E-19 0.		0.
4050.00-4100.00	3.98E+13 0.		2.91E-23 0.		0.	6.59E-19 0.		0.
4100.00-4150.00	3.98E+13 0.		3.14E-23 0.		0.	6.03E-19 0.		0.
4150.00-4200.00	4.04E+13 0.		3.99E-23 0.		0.	6.24E-19 0.		0.
4200.00-4250.00	4.02E+13 0.		6.54E-23 0.		0.	0.		0.
4250.00-4300.00	3.88E+13 0.		6.83E-23 0.		0.	0.		0.
4300.00-4350.00	3.96E+13 0.		8.66E-23 0.		0.	0.		0.
4350.00-4400.00	4.50E+13 0.		1.25E-22 0.		0.	0.		0.
4400.00-4450.00	4.78E+13 0.		1.49E-22 0.		0.	0.		0.
4450.00-4500.00	4.96E+13 0.		1.71E-22 0.		0.	0.		0.
4500.00-4550.00	4.98E+13 0.		2.12E-22 0.		0.	0.	6.00E-20	0.
4550.00-4600.00	4.96E+13 0.		3.57E-22 0.		0.	0.	1.00E-19	0.
4600.00-4650.00	5.00E+13 0.		3.68E-22 0.		0.	0.	8.00E-20	0.
4650.00-4700.00	5.10E+13 0.		4.06E-22 0.		0.	0.	1.20E-19	0.
4700.00-4750.00	5.22E+13 0.		4.89E-22 0.		0.	0.	1.80E-19	0.
4750.00-4800.00	5.18E+13 0.		7.11E-22 0.		0.	0.	2.40E-19	0.
4800.00-4850.00	4.92E+13 0.		8.43E-22 0.		0.	0.	2.10E-19	0.
4850.00-4900.00	4.88E+13 0.		8.28E-22 0.		0.	0.	2.40E-19	0.
4900.00-4950.00	5.06E+13 0.		9.09E-22 0.		0.	0.	2.40E-19	0.
4950.00-5000.00	4.96E+13 0.		1.22E-21 0.		0.	0.	3.00E-19	0.
5000.00-5050.00	4.98E+13 0.		1.62E-21 0.		0.	0.	3.00E-19	0.
5050.00-5100.00	5.00E+13 0.		1.58E-21 0.		0.	0.	4.00E-19	0.
5100.00-5150.00	4.86E+13 0.		1.60E-21 0.		0.	0.	4.50E-19	0.
5150.00-5200.00	4.86E+13 0.		1.78E-21 0.		0.	0.	5.00E-19	0.
5200.00-5250.00	5.04E+13 0.		2.07E-21 0.		0.	0.	5.00E-19	0.
5250.00-5300.00	5.16E+13 0.		2.55E-21 0.		0.	0.	6.00E-19	0.
5300.00-5350.00	5.28E+13 0.		2.74E-21 0.		0.	0.	6.00E-19	0.
5350.00-5400.00	5.34E+13 0.		2.88E-21 0.		0.	0.	5.00E-19	0.
5400.00-5450.00	5.40E+13 0.		3.07E-21 0.		0.	0.	4.00E-19	0.
5450.00-5500.00	5.36E+13 0.		3.17E-21 0.		0.	0.	6.00E-19	0.
5500.00-5550.00	5.32E+13 0.		3.36E-21 0.		0.	0.	6.50E-19	0.
5550.00-5600.00	5.34E+13 0.		3.88E-21 0.		0.	0.	7.00E-19	0.
5600.00-5650.00	5.34E+13 0.		4.31E-21 0.		0.	0.	6.00E-19	0.
5650.00-5700.00	5.34E+13 0.		4.67E-21 0.		0.	0.	6.00E-19	0.

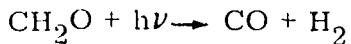
Table B1. Solar Fluxes and Absorption Cross Sections for O₂, O₃, H₂O, H₂O₂, NO₂, NO₃, and N₂O (Contd)

WAVELENGTH BAND ANGSTROMS	FLUX PH/CM ² /S/A	CROSS SECTIONS (CM ²)						
		O2	O3	H2O	H2O2	NO2	NO3	N2O
1792.60-1803.60	1.01E+10	0.	7.77E-19	1.02E-18	0.	0.	0.	0.
1803.60-1816.40	1.27E+10	0.	7.60E-19	6.98E-19	0.	0.	0.	0.
1816.40-1830.60	1.42E+10	0.	7.33E-19	3.64E-19	0.	0.	0.	0.
1830.60-1846.20	1.29E+10	0.	6.99E-19	1.80E-19	0.	0.	0.	0.
1846.20-1863.40	1.31E+10	0.	6.56E-19	6.51E-20	0.	0.	0.	0.
1863.40-1882.20	1.63E+10	0.	6.06E-19	0.	0.	0.	0.	0.
1882.20-1902.40	1.92E+10	0.	5.45E-19	0.	9.79E-19	0.	0.	0.
1902.40-1924.00	2.29E+10	0.	4.86E-19	0.	8.10E-19	5.25E-19	0.	1.26E-19
1924.00-1947.00	3.61E+10	0.	4.36E-19	0.	7.03E-19	4.83E-19	0.	9.00E-20
1947.00-1971.80	4.68E+10	0.	4.01E-19	0.	6.15E-19	4.00E-19	0.	8.00E-20
1971.80-1985.00	6.21E+10	0.	3.42E-19	0.	5.64E-19	3.50E-19	0.	7.00E-20
1985.00-2000.00	6.85E+10	0.	3.28E-19	0.	5.43E-19	3.30E-19	0.	5.50E-20
2000.00-2025.00	7.33E+10	0.	3.26E-19	0.	5.16E-19	3.50E-19	0.	4.00E-20
2025.00-2050.00	8.74E+10	0.	3.28E-19	0.	4.79E-19	4.00E-19	0.	2.50E-20
2050.00-2061.86	1.10E+11	1.05E-23	3.51E-19	0.	4.66E-19	4.20E-19	0.	2.00E-20
2061.86-2083.33	1.96E+11	1.00E-23	4.11E-19	0.	4.42E-19	4.30E-19	0.	1.70E-20
2083.33-2105.26	3.33E+11	9.55E-24	4.84E-19	0.	4.22E-19	4.40E-19	0.	1.10E-20
2105.26-2127.66	4.21E+11	8.93E-24	6.26E-19	0.	3.92E-19	4.54E-19	0.	4.00E-21
2127.66-2150.54	4.63E+11	8.28E-24	8.57E-19	0.	3.63E-19	4.58E-19	0.	2.00E-21
2150.54-2173.91	5.73E+11	7.60E-24	1.17E-18	0.	3.40E-19	4.62E-19	0.	1.10E-21
2173.91-2197.80	5.53E+11	6.92E-24	1.52E-18	0.	3.19E-19	4.50E-19	0.	1.00E-21
2197.80-2222.22	7.08E+11	6.28E-24	1.97E-18	0.	2.98E-19	4.25E-19	0.	4.00E-22
2222.22-2247.19	7.21E+11	5.65E-24	2.55E-18	0.	2.67E-19	4.00E-19	0.	2.80E-22
2247.19-2272.73	7.13E+11	5.03E-24	3.24E-18	0.	2.49E-19	3.50E-19	0.	2.00E-22
2272.73-2298.85	8.65E+11	4.40E-24	4.00E-18	0.	2.36E-19	3.00E-19	0.	1.20E-22
2298.85-2325.58	8.98E+11	3.76E-24	4.83E-18	0.	2.16E-19	2.50E-19	0.	7.00E-23
2325.58-2352.94	8.22E+11	3.09E-24	5.79E-18	0.	1.99E-19	2.00E-19	0.	5.00E-23
2352.94-2380.95	7.89E+11	2.44E-24	6.86E-18	0.	1.81E-19	1.18E-19	0.	3.00E-23
2380.95-2409.64	8.09E+11	1.75E-24	7.97E-18	0.	1.66E-19	1.00E-19	0.	2.00E-23
2409.64-2439.02	8.51E+11	6.74E-25	9.00E-18	0.	1.48E-19	7.00E-20	0.	1.10E-23
2439.02-2469.14	9.07E+11	0.	1.00E-17	0.	1.30E-19	5.00E-20	0.	9.00E-24
2469.14-2500.00	9.33E+11	0.	1.07E-17	0.	1.22E-19	3.60E-20	0.	8.00E-24
2500.00-2531.65	9.54E+11	0.	1.11E-17	0.	1.08E-19	2.50E-20	0.	7.00E-24
2531.65-2564.10	1.22E+12	0.	1.12E-17	0.	9.30E-20	1.50E-20	0.	6.00E-24
2554.10-2597.40	2.14E+12	0.	1.11E-17	0.	8.43E-20	1.70E-20	0.	5.00E-24
2597.40-2631.58	1.28E+12	0.	1.03E-17	0.	7.49E-20	1.90E-20	0.	5.00E-24
2631.58-2666.67	3.19E+12	0.	9.43E-18	0.	6.39E-20	2.10E-20	0.	5.00E-24
2666.67-2702.70	3.47E+12	0.	8.23E-18	0.	5.64E-20	2.90E-20	0.	5.00E-24
2702.70-2739.73	3.13E+12	0.	6.81E-18	0.	5.00E-20	3.70E-20	0.	5.20E-24
2739.73-2777.78	3.13E+12	0.	5.31E-18	0.	4.33E-20	4.20E-20	0.	5.90E-24
2777.78-2816.90	3.53E+12	0.	3.99E-18	0.	3.62E-20	5.00E-20	0.	6.00E-24
2816.90-2857.14	4.22E+12	0.	2.84E-18	0.	3.27E-20	6.50E-20	0.	5.40E-24
2857.14-2898.55	5.94E+12	0.	1.92E-18	0.	2.91E-20	8.00E-20	0.	5.20E-24
2898.55-2941.18	9.15E+12	0.	1.14E-18	0.	2.53E-20	9.00E-20	0.	5.00E-24
2941.18-2985.07	9.09E+12	0.	6.60E-19	0.	2.15E-20	1.20E-19	0.	5.10E-24
2985.07-3030.30	6.53E+12	0.	3.69E-19	0.	1.75E-20	1.33E-19	0.	5.00E-24
3030.30-3076.92	1.09E+13	0.	1.97E-19	0.	0.	1.65E-19	0.	4.20E-24
3076.92-3110.00	2.57E+13	0.	1.05E-19	0.	0.	1.97E-19	0.	3.40E-24
3110.00-3140.00	1.21E+13	0.	5.23E-20	0.	0.	2.24E-19	0.	1.70E-24
3140.00-3200.00	1.39E+13	0.	2.91E-20	0.	0.	2.61E-19	0.	0.

Table B1. Solar Fluxes and Absorption Cross Sections for O₂, O₃, H₂O,
H₂O₂, NO₂, NO₃, and N₂O

WAVELENGTH BAND ANGSTROMS	FLUX PH/CM ² /S/A	CROSS SECTIONS (CM ²)						
		O2	O3	H2O	H2O2	NO2	NO3	N2O
1215.67-1215.67	3.00E+11	1.00E-20	2.32E-17	0.	0.	0.	0.	0.
1162.79-1169.59	1.51E+07	2.00E-20	7.80E-18	0.	0.	0.	0.	0.
1169.59-1176.47	3.87E+07	1.25E-18	7.97E-18	0.	0.	0.	0.	0.
1176.47-1183.43	1.61E+07	2.55E-19	8.66E-18	0.	0.	0.	0.	0.
1183.43-1190.48	1.76E+07	3.00E-20	9.51E-18	0.	0.	0.	0.	0.
1190.48-1197.60	2.55E+07	3.75E-19	1.25E-17	0.	0.	0.	0.	0.
1197.60-1204.82	2.87E+07	4.45E-18	1.84E-17	0.	0.	0.	0.	0.
1204.82-1212.12	1.01E+08	8.35E-18	2.19E-17	0.	0.	0.	0.	0.
1212.12-1219.51	3.08E+08	6.00E-19	2.30E-17	0.	0.	0.	0.	0.
1219.51-1226.99	4.91E+08	2.35E-19	2.26E-17	1.99E-17	0.	0.	0.	0.
1226.99-1234.57	1.80E+08	4.50E-19	2.06E-17	6.15E-18	0.	0.	0.	0.
1234.57-1242.24	1.51E+08	3.35E-19	1.30E-17	5.16E-18	0.	0.	0.	0.
1242.24-1250.00	9.69E+07	1.75E-17	8.91E-18	1.10E-17	0.	0.	0.	0.
1250.00-1257.86	1.05E+08	8.95E-19	7.24E-18	6.39E-18	0.	0.	0.	0.
1257.86-1265.82	1.19E+08	4.30E-19	6.09E-18	7.13E-18	0.	0.	0.	0.
1265.82-1273.89	1.03E+08	1.10E-19	5.66E-18	7.84E-18	0.	0.	0.	0.
1273.89-1282.05	7.96E+07	2.05E-19	5.87E-18	7.48E-18	0.	0.	0.	0.
1282.05-1290.32	5.71E+07	4.43E-19	6.47E-18	7.82E-18	0.	0.	0.	0.
1290.32-1298.70	7.16E+07	5.55E-19	8.14E-18	7.64E-18	0.	0.	0.	0.
1298.70-1307.19	3.89E+08	4.20E-19	1.24E-17	6.71E-18	0.	0.	0.	0.
1307.19-1315.79	2.06E+08	6.85E-19	1.52E-17	6.75E-18	0.	0.	0.	0.
1315.79-1324.50	8.68E+07	1.45E-18	1.47E-17	5.67E-18	0.	0.	0.	0.
1324.50-1333.33	1.73E+08	2.25E-18	1.51E-17	4.93E-18	0.	0.	0.	0.
1333.33-1342.28	2.79E+08	2.30E-18	1.51E-17	5.00E-18	0.	0.	0.	0.
1342.28-1351.35	1.03E+08	4.55E-18	1.65E-17	3.69E-18	0.	0.	0.	0.
1351.35-1360.54	1.62E+08	7.23E-18	1.54E-17	3.36E-18	0.	0.	0.	0.
1360.54-1369.86	1.23E+08	9.50E-18	1.35E-17	2.42E-18	0.	0.	0.	0.
1369.86-1379.31	1.37E+08	1.23E-17	1.05E-17	1.80E-18	0.	0.	0.	0.
1379.31-1388.89	1.20E+08	1.32E-17	7.97E-18	1.60E-18	0.	0.	0.	0.
1388.89-1408.45	2.93E+08	1.36E-17	7.17E-18	1.19E-18	0.	0.	0.	0.
1408.45-1428.57	2.10E+08	1.40E-17	6.28E-18	7.00E-19	0.	0.	0.	0.
1428.57-1449.28	2.49E+08	1.48E-17	5.66E-18	5.03E-19	0.	0.	0.	0.
1449.28-1470.59	2.94E+08	1.41E-17	5.23E-18	4.96E-19	0.	0.	0.	0.
1470.59-1492.54	4.06E+08	1.29E-17	4.47E-18	6.87E-19	0.	0.	0.	0.
1492.54-1515.15	4.55E+08	1.15E-17	3.69E-18	9.45E-19	0.	0.	0.	0.
1515.15-1538.46	5.96E+08	9.91E-18	2.93E-18	1.38E-18	0.	0.	0.	0.
1538.46-1562.50	9.36E+08	8.24E-18	2.19E-18	1.84E-18	0.	0.	0.	0.
1562.50-1587.30	8.59E+09	6.58E-18	1.63E-18	2.44E-18	0.	0.	0.	0.
1587.30-1612.90	8.05E+08	4.97E-18	1.20E-18	3.20E-18	0.	0.	0.	0.
1612.90-1639.34	1.29E+09	3.45E-18	9.77E-19	3.91E-18	0.	0.	0.	0.
1639.34-1665.67	2.06E+09	2.08E-18	8.66E-19	4.64E-18	0.	0.	0.	0.
1665.67-1694.92	2.84E+09	1.23E-18	8.14E-19	4.57E-18	0.	0.	0.	0.
1694.92-1724.14	4.72E+09	7.22E-19	8.17E-19	4.35E-18	0.	0.	0.	0.
1724.14-1759.13	4.93E+09	4.56E-19	8.57E-19	3.57E-18	0.	0.	0.	0.
1759.13-1780.00	5.49E+09	2.74E-19	8.40E-19	2.64E-18	0.	0.	0.	0.
1780.00-1782.20	6.25E+09	0.	8.18E-19	2.64E-18	0.	0.	0.	0.
1783.20-1789.60	7.60E+09	0.	8.04E-19	2.22E-18	0.	0.	0.	0.
1789.60-1794.60	9.14E+09	0.	8.00E-19	2.04E-18	0.	0.	0.	0.
1794.60-1798.60	8.42E+09	0.	7.75E-19	1.83E-18	0.	0.	0.	0.
1798.60-1802.60	8.06E+09	0.	7.89E-19	1.42E-18	0.	0.	0.	0.

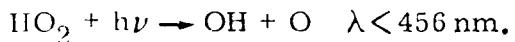
and



for wavelengths from 240 to 360 nm. The absorption cross sections for this molecule are taken from McQuigg and Calvert^{B14} and Calvert et al.^{B15}. These authors also give the quantum efficiencies for these processes. Table B19 lists the total photodissociation rate coefficients for formaldehyde. If both of the above processes are considered to occur for this molecule, then the values presented in Table B19 must be multiplied by the relative efficiencies as given by McQuigg and Calvert^{B14} and Calvert et al^{B15} for each of these reactions.

B14. THE HYDROPEROXYL RADICAL

The primary photodissociation process for the hydroperoxyl radical is



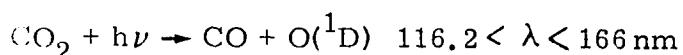
The absorption cross sections used for HO₂ are a composite of the measurements of Hochanadel et al^{B16} and Paukert and Johnston.^{B17} The photodissociation rate coefficients for this radical are listed in Table B20.

-
- B14. McQuigg, R.D., and Calvert, J.G. (1969) The photodecomposition of CH₂O, CD₂O, CHDO and CH₂O-CD₂O mixtures at xenon flash lamp intensities, J. Am. Chem. Soc. 91:1590-1599.
- B15. Calvert, J.G., Kerr, J.A., Demerjian, K.L., and McQuigg, R.D. (1972) Photolysis of formaldehyde as a hydrogen atom source in the lower atmosphere, Science 175:751-752.
- B16. Hochanadel, C.J., Ghormley, A., and Ogren, P.J. (1972) Absorption spectrum and reaction kinetics of the HO₂ radical in the gas phase, J. Chem. Phys. 56:4426-4432.
- B17. Paukert, T.T., and Johnston, A.S. (1972) Spectra and kinetics of the hydroperoxyl free radical in the gas phase, J. Chem. Phys. 56:2824-2833.

for wavelengths from 116 to 164 nm. Most of the CH_4 photodissociation occurs at the Lyman α line 121.6 nm. The values listed in Table B16 are the total photodissociation rate coefficients for methane. If the above processes have equal efficiencies, then each process would have rate coefficients equal to half of the values shown in the table. The absorption cross sections for this molecule are taken from Watanabe et al.^{B10}

B12. CARBON DIOXIDE

The photolysis of carbon dioxide can be proceed by two different processes:



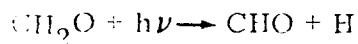
and



The absorption cross sections for this molecule are taken from Inn et al,^{B11} Heimerl,^{B12} and Ogawa.^{B13} The photodissociation rate coefficients for the processes yielding excited atomic oxygen are listed in Table B17, and those for the processes yielding ground state atomic oxygen are listed in Table B18.

B13. FORMALDEHYDE

Although formaldehyde can be photodissociated in several ways, the primary dissociative processes are

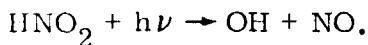


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- (1) Watanabe, K., Zelikoff, M., and Inn, E. C. Y. (1953) Absorption Coefficient of Several Atmospheric Gases, Geophys. Res. Paper, 21, Tech Rpt. AFCRL-53-23.
 - (2) Inn, E. C. Y., Watanabe, K., and Zelikoff, M. (1953) Absorption coefficients of gases in the vacuum ultraviolet. Part III. CO_2 , J. Chem. Phys. 21:1648-1651.
 - (3) Heimerl, J. (1970) CO_2 absorption coefficient 1655-1825 A, J. Geophys. Res. 75:5574-5575.
 - (4) Ogawa, M. (1971) Absorption cross sections of O_2 and CO_2 continua in the Schumann and UV regions, J. Chem. Phys. 54:2550-2556.

cule are taken from Jones and Wulf^{B9} and Johnston and Graham.¹³ Its photodissociation rate coefficients are listed in Table B13.

B9. NITROUS ACID

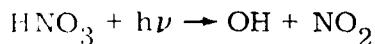
The photodissociation of nitrous acid probably proceeds as



The absorption cross sections for this molecule between 300 and 390 nm are those given by Johnston and Graham.¹³ Since O₂ is not an efficient absorber of radiation in this wavelength interval, the photodissociation rate coefficients for HNO₂ are listed in Table B14 only as a function of the column density of ozone.

B10. NITRIC ACID

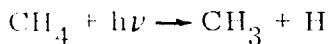
The primary photolytic process for nitric acid is



over the broad wavelength region from 192 to 325 nm. The values for the absorption cross sections for this gas are taken from Johnston and Graham.¹³ The nitric acid photodissociation rate coefficients are listed in Table B15.

B11. METHANE

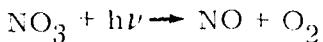
The photodissociation processes for methane are



and



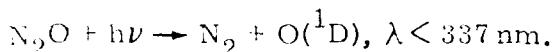
B9. Jones, E.J., and Wulf, O.R. (1937) The absorption coefficient of nitrogen pentoxide in the ultraviolet and the visible absorption spectrum of NO₃, *J. Chem. Phys.* 5:873-877.



for wavelengths longer than 578 nm. For wavelengths below 578 nm, it is not known whether the products are $\text{NO}_2 + \text{O}$ or $\text{NO} + \text{O}_2$. In Table B11, all the photodissociation of this radical is assumed to produce $\text{NO} + \text{O}_2$. The absorption cross sections for this molecule are taken from Johnston and Graham.¹³ Since molecular oxygen is not an efficient absorber for radiations above 450 nm, the photodissociation rate coefficients for this molecule are given only as a function of the ozone column density.

B7. NITROUS OXIDE

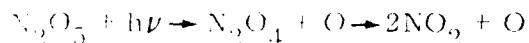
The primary photolytic process for nitrous oxide according to Preston and Baer^{B6} is



Below 210 nm, oxygen atoms in the ^1S state may be produced. For the calculations presented in this report, Table B12 was computed with all of the photodissociation of N_2O producing $\text{O}(\text{^1D})$. The absorption cross sections below 205 nm are from Zelikoff et al.^{B7} and above 205 nm, from Bates and Hayes.^{B8}

B8. DINITROGEN PENTOXIDE

The primary processes in the photolysis of N_2O_5 are not well understood. Jones and Wulf^{B9} attribute the photodissociation of this molecule to



at wavelengths from 210 to 380 nm. The absorption cross sections for this mol-

-
- (a) Preston, R.F., and Baer, R.F. (1971) Primary processes in the photolysis of nitrous oxide, *J. Chem. Phys.*, 54, 3347-3348.
 - (b) Zelikoff, M., Watanabe, K., and Imamura, E.C.Y. (1953) Absorption coefficients of gases in the vacuum ultraviolet. Part II. Nitrous oxide, *J. Chem. Phys.*, 21, 1643-1647.
 - (c) Bates, D.R., and Hays, P.B. (1967) Atmospheric nitrous oxide, *Planet. Space Sci.*, 15, 161-177.

Table B3. Total photodissociation rate coefficients for the process $O_2 + hv \rightarrow O + O(^1D)$ (Contd)

	O3 COLUMN DENSITY				
	5.0E+18	1.0E+19	2.0E+19	5.0E+19	1.0E+20
1.0E+13	6.79854E-09	9.47067E-11	2.23187E-14	4.24957E-25	7.77344E-43
2.0E+13	6.79822E-09	9.47055E-11	2.23184E-14	4.24953E-25	7.77336E-43
5.0E+13	6.79788E-09	9.47016E-11	2.23177E-14	4.24940E-25	7.77312E-43
1.0E+14	6.79731E-09	9.46952E-11	2.23164E-14	4.24919E-25	7.77273E-43
2.0E+14	6.79618E-09	9.46824E-11	2.23139E-14	4.24876E-25	7.77195E-43
5.0E+14	6.79277E-09	9.46440E-11	2.23062E-14	4.24746E-25	7.76959E-43
1.0E+15	6.78709E-09	9.45800E-11	2.22935E-14	4.24531E-25	7.76567E-43
2.0E+15	6.77575E-09	9.44522E-11	2.22680E-14	4.24101E-25	7.75783E-43
5.0E+15	6.74192E-09	9.40702E-11	2.21919E-14	4.22813E-25	7.73437E-43
1.0E+16	6.68610E-09	9.34384E-11	2.20657E-14	4.20676E-25	7.69544E-43
2.0E+16	6.57656E-09	9.21926E-11	2.18161E-14	4.16438E-25	7.61821E-43
5.0E+16	6.26383E-09	8.85918E-11	2.10884E-14	4.04011E-25	7.39145E-43
1.0E+17	5.79071E-09	8.30146E-11	2.19426E-14	3.84213E-25	7.02941E-43
2.0E+17	4.99391E-09	7.32432E-11	1.78774E-14	3.47802E-25	6.36107E-43
5.0E+17	3.39109E-09	5.19651E-11	1.31032E-14	2.59836E-25	4.73406E-43
1.0E+18	2.01269E-09	3.17454E-11	8.18451E-15	1.63393E-25	2.93529E-43
2.0E+18	8.94428E-10	1.41535E-11	3.62480E-15	1.8981E-26	1.09807E-43
5.0E+18	1.58965E-10	2.41572E-12	5.68530E-16	8.57694E-27	4.23429E-43
1.0E+19	2.18886E-11	3.24585E-13	7.18573E-17	8.26195E-28	6.12377E-46
2.0E+19	1.07208E-12	1.60175E-14	3.57869E-18	4.01515E-29	2.29144E-47
5.0E+19	2.75050E-16	4.12506E-18	9.27565E-22	1.05462E-32	6.06327E-51
1.0E+20	3.08714E-22	4.62933E-24	1.04099E-27	1.18365E-38	6.80543E-57
2.0E+20	4.49203E-31	5.83243E-36	1.31152E-39	1.49127E-50	8.57407E-69
5.0E+20	1.08004E-33	1.24725E-50	2.62283E-75	2.98228E-86	1.71467E-104
1.0E+21	4.90335E-38	5.66250E-55	7.55159E-89	9.46683E-146	5.66818E-141
2.0E+21	1.01066E-48	1.16713E-63	1.55650E-97	3.69182E-199	1.79928E-200
5.0E+21	8.84982E-73	1.0220E-89	1.36295E-123	3.23274E-225	0.
1.0E+22	3.29220E-116	3.80190E-133	5.07028E-167	1.20260E-268	0.

Table B5. Total Photodissociation Rate Coefficients for the Process $O_2 + h\nu \rightarrow O + O$ for Wavelengths From 175 to 242.2 nm

	O3 COLUMN DENSITY									
	1.0E+13	2.0E+13	5.0E+13	1.0E+14	2.0E+14	5.0E+14	1.0E+15	2.0E+15	5.0E+15	1.0E+16
1.0E+17	8.36791E-08	8.36784E-08	8.36764E-08	8.36731E-08	8.36664E-08	8.36464E-08	8.36131E-08	8.35466E-08	8.33473E-08	8.30164E-08
2.0E+17	7.96609E-08	7.96602E-08	7.96583E-08	7.96551E-08	7.96488E-08	7.96290E-08	7.95980E-08	7.95346E-08	7.93448E-08	7.90296E-08
5.0E+17	6.94462E-08	6.94456E-08	6.94440E-08	6.94412E-08	6.94357E-08	6.94190E-08	6.93913E-08	6.93359E-08	6.91701E-08	6.88949E-08
1.0E+18	5.78378E-08	5.78373E-08	5.78359E-08	5.78336E-08	5.78290E-08	5.78151E-08	5.77919E-08	5.77456E-08	5.76070E-08	5.73770E-08
2.0E+18	4.48169E-08	4.48165E-08	4.48136E-08	4.48113E-08	4.48100E-08	4.47991E-08	4.47811E-08	4.47450E-08	4.46369E-08	4.44575E-08
5.0E+18	2.84480E-08	2.84478E-08	2.84471E-08	2.84459E-08	2.84435E-08	2.84365E-08	2.84248E-08	2.84014E-08	2.83314E-08	2.82154E-08
1.0E+19	1.87404E-08	1.87402E-08	1.87390E-08	1.87374E-08	1.87357E-08	1.87326E-08	1.87246E-08	1.87087E-08	1.86612E-08	1.85824E-08
2.0E+19	1.19434E-08	1.19433E-08	1.19429E-08	1.19424E-08	1.19413E-08	1.19381E-08	1.19327E-08	1.19220E-08	1.18898E-08	1.18367E-08
5.0E+19	6.78508E-09	6.78501E-09	6.78496E-09	6.78495E-09	6.78375E-09	6.78166E-09	6.78119E-09	6.77119E-09	6.75038E-09	6.71603E-09
1.0E+20	4.62169E-09	4.62163E-09	4.62147E-09	4.62120E-09	4.62065E-09	4.61901E-09	4.61628E-09	4.61083E-09	4.59459E-09	4.56784E-09
2.0E+20	3.25814E-09	3.25810E-09	3.25796E-09	3.25773E-09	3.25728E-09	3.25592E-09	3.25365E-09	3.24914E-09	3.23568E-09	3.21355E-09
5.0E+20	3.25604E-09	3.25599E-09	3.25586E-09	3.25563E-09	3.25518E-09	3.25382E-09	3.25155E-09	3.24704E-09	3.23359E-09	3.21149E-09
1.0E+21	2.19371E-09	2.19963E-09	2.19956E-09	2.19937E-09	2.19898E-C9	2.19763E-09	2.19591E-09	2.19209E-09	2.18071E-09	2.16203E-09
2.0E+21	1.75574E-09	1.75574E-09	1.75563E-09	1.75561E-09	1.75546E-09	1.75540E-09	1.75525E-09	1.74870E-09	1.7385Y-09	1.72086E-09
5.0E+21	1.49053E-09	1.49050E-09	1.49040E-09	1.49023E-09	1.48689E-C9	1.48689E-09	1.48687E-09	1.48381E-09	1.47379E-09	1.45738E-09
1.0E+22	1.29543E-09	1.29540E-09	1.29530E-09	1.29514E-09	1.29482E-09	1.29385E-09	1.29223E-09	1.28901E-09	1.27944E-09	1.26378E-09
2.0E+22	1.16775E-09	1.16775E-09	1.16766E-09	1.16750E-09	1.16720E-09	1.16627E-09	1.16474E-09	1.16168E-09	1.15259E-09	1.13770E-09
5.0E+22	9.58054E-10	9.580627E-10	9.58046E-10	9.58010E-10	9.580139E-10	9.580128E-10	9.580077E-10	9.53288E-C10	9.45287E-10	9.32204E-10
1.0E+23	7.03317E-10	7.03295E-10	7.03229E-10	7.03118E-10	7.02897E-10	7.02234E-10	7.01131E-10	6.98933E-10	6.92405E-10	6.81137E-10
2.0E+23	4.12219E-10	4.12204E-10	4.12158E-10	4.12081E-10	4.11928E-10	4.11471E-10	4.10709E-10	4.09192E-10	4.04690E-10	3.97346E-10
5.0E+23	1.18661E-10	1.18655E-10	1.18636E-10	1.18606E-10	1.18545E-10	1.18362E-10	1.18058E-10	1.17454E-10	1.15664E-10	1.12756E-10
1.0E+24	3.05886E-11	3.05666E-11	3.05601E-11	3.05702E-11	3.05498E-11	3.04888E-11	3.03869E-11	3.01848E-11	2.95873E-11	2.86213E-11
2.0E+24	8.54771E-12	8.54716E-12	8.54551E-12	8.54275E-12	8.53727E-12	8.52032E-12	8.49249E-12	8.43918E-12	8.27895E-12	8.02063E-12
5.0E+24	2.90922E-12	2.80815E-12	2.90794E-12	2.90758E-12	2.90473E-12	2.90117E-12	2.89410E-12	2.88924E-12	2.83718E-12	2.33181E-12
1.0E+25	2.34202E-12	2.34200E-12	2.34193E-12	2.34193E-12	2.34163E-12	2.34163E-12	2.33794E-12	2.33794E-12	2.32165E-12	2.32120E-12
2.0E+25	2.32211E-12	2.32209E-12	2.32204E-12	2.32194E-12	2.32176E-12	2.32120E-12	2.31839E-12	2.31270E-12	2.30346E-12	

Table 11. Total Uncondensatation Rate Coefficients for the Process $O_2 + h\nu \rightarrow O + O$ for Wavelengths From 175 to 242.2 nm (Cont.)

O3 COLUMN DENSITY									
	2.0E+16	5.0E+16	1.0E+17	2.0E+17	5.0E+17	1.0E+18	2.0E+18	5.0E+18	1.0E+19
1.0E+17	8.23593E-08	8.04246E-08	7.73159E-08	7.14945E-08	5.66739E-08	3.86203E-08	1.80502E-08	1.93497E-09	6.22227E-11
2.0E+17	7.84037E-08	7.6562E-08	7.36013E-08	6.80506E-08	5.39608E-08	3.67881E-08	1.72141E-08	1.85505E-09	6.05895E-11
5.0E+17	6.83484E-08	6.67403E-08	6.41590E-08	5.93332E-08	4.70661E-08	3.21332E-08	1.50907E-08	1.65214E-09	5.64206E-11
1.0E+18	5.69205E-08	5.55781E-08	5.341E-08	4.94087E-08	3.92220E-08	2.68312E-08	1.26659E-08	1.41812E-09	5.14674E-11
2.0E+18	4.41017E-08	4.30567E-08	4.13851E-08	3.82738E-08	3.08730E-08	2.04156E-08	9.93541E-09	1.15254E-09	4.57406E-11
5.0E+18	2.79855E-08	2.73124E-08	2.62414E-08	2.42635E-08	1.93198E-08	1.33475E-08	6.46744E-09	8.07455E-10	3.77590E-11
1.0E+19	1.84265E-08	1.79722E-08	1.72547E-08	1.59439E-08	1.27178E-08	8.85386E-09	4.38021E-09	5.93023E-10	3.23310E-11
2.0E+19	1.17318E-08	1.14281E-08	1.09541E-08	1.01032E-08	8.63595E-09	5.66293E-09	2.87414E-09	4.28336E-10	2.73964E-11
5.0E+19	6.64658E-08	6.45547E-08	6.06434E-08	5.64812E-09	4.48576E-09	3.17666E-09	1.66625E-09	2.81839E-10	2.188888E-11
1.0E+20	4.51555E-09	4.36707E-09	4.14439E-09	3.76924E-09	2.96396E-09	2.10352E-09	1.12840E-09	2.09148E-10	1.84843E-11
2.0E+20	2.35411E-09	2.04920E-09	2.07263E-09	2.057918E-09	1.99268E-09	1.40941E-09	7.70327E-10	1.55452E-10	1.53759E-11
5.0E+20	3.16638E-09	3.04728E-09	2.86888E-09	2.57769E-09	2.199168E-09	1.40879E-09	7.70033E-10	1.55397E-10	1.53705E-11
1.0E+21	2.12576E-09	2.02486E-09	1.87920E-09	1.65023E-09	1.22740E-09	8.53232E-10	4.72751E-10	1.04948E-10	1.17558E-11
2.0E+21	1.66732E-09	1.59453E-09	1.46199E-09	1.25738E-09	9.00217E-10	6.10129E-10	3.36480E-10	7.83265E-11	9.43953E-12
5.0E+21	1.42560E-09	1.33794E-09	1.21157E-09	1.02460E-09	7.05255E-10	4.63424E-10	2.51417E-10	5.99429E-11	7.63448E-12
1.0E+22	1.23347E-09	1.15007E-09	1.03236E-09	8.55318E-10	5.64536E-10	3.56948E-10	1.88097E-10	4.50794E-11	6.02269E-12
2.0E+22	1.10893E-09	1.02987E-09	9.18651E-10	7.52459E-10	4.64043E-10	2.98625E-10	1.54032E-10	3.66700E-11	5.00513E-12
5.0E+22	9.06934E-10	8.37738E-10	7.40991E-10	5.98057E-10	3.73317E-10	2.24810E-10	1.14036E-10	2.71906E-11	3.81188E-12
1.0E+23	6.61173E-10	6.05122E-10	5.27515E-10	4.14905E-10	2.45244E-10	1.40952E-10	6.87819E-11	1.59798E-11	2.25926E-12
2.0E+23	3.83235E-10	3.45100E-10	2.93210E-10	2.20309E-10	1.18551E-10	6.35160E-11	2.96893E-11	6.85088E-12	1.00007E-12
5.0E+23	1.07213E-10	9.25449E-11	7.34312E-11	4.85877E-11	2.03970E-11	1.964924E-12	4.57852E-12	1.21106E-12	2.05296E-13
1.0E+24	2.67951E-11	2.20719E-11	1.62007E-11	9.26161E-12	2.93456E-12	1.29966E-12	5.41249E-13	6.31005E-14	4.84552E-15
2.0E+24	7.53517E-12	6.29823E-12	4.80552E-12	3.12606E-12	1.65148E-12	1.03930E-12	4.63488E-13	4.24045E-14	1.07364E-15
5.0E+24	2.77466E-12	2.60425E-12	2.38066E-12	2.07309E-12	1.55767E-12	1.03654E-12	4.63449E-13	4.24042E-14	1.07363E-15
1.0E+25	2.30156E-12	2.24295E-12	2.15012E-12	1.97922E-12	1.55133E-12	1.03647E-12	4.63449E-13	4.24042E-14	1.07363E-15
2.0E+25	2.28494E-12	2.23026E-12	2.14203E-12	1.97593E-12	1.55111E-12	1.03647E-12	4.63449E-13	4.24042E-14	1.07363E-15

O2 COLUMN DENSITY

Table B5. Total Photodissociation Rate Coefficients for the Process $O_2 + h\nu \rightarrow O + O$ for Wavelengths From 175 to 242.2 nm (Contd)

	O3 COLUMN DENSITY				
	5.0E+19	1.0E+20	2.0E+20	5.0E+20	1.0E+21
1.0E+17	5.73391E-18	5.25966E-25	2.73918E-39	6.48784E-82	7.833393-153
2.0E+17	8.73401E-18	5.25969E-25	2.73918E-39	6.48784E-82	7.833393-153
5.0E+17	8.73431E-18	5.25972E-25	2.73915E-39	6.48771E-82	7.833377-153
1.0E+18	8.70961E-18	5.25049E-25	2.73786E-39	6.48769E-82	7.833385-153
2.0E+18	8.70936E-18	5.25076E-25	2.73786E-39	6.48763E-82	7.833385-153
5.0E+18	8.70805E-18	5.25147E-25	2.73783E-39	6.48732E-82	7.833366-153
1.0E+19	8.68652E-18	5.24339E-25	2.73649E-39	6.48685E-82	7.833351-153
2.0E+19	8.64264E-18	5.22092E-25	2.72619E-39	6.46929E-82	7.82302-153
5.0E+19	8.55010E-18	5.19131E-25	2.72091E-39	6.46673E-82	7.82211-153
1.0E+20	8.43410E-18	5.15784E-25	2.71819E-39	6.47105E-82	7.82488-153
2.0E+20	8.21772E-18	5.09124E-25	2.70931E-39	6.47191E-82	7.82630-153
5.0E+20	8.21654E-18	5.09100E-25	2.70929E-39	6.47191E-82	7.82630-153
1.0E+21	7.75047E-18	4.94050E-25	2.68714E-39	6.46964E-82	7.82824-153
2.0E+21	7.29024E-18	4.77961E-25	2.65499E-39	6.44707E-82	7.81927-153
5.0E+21	6.81167E-18	4.59358E-25	2.60309E-39	6.36972E-82	7.73855-153
1.0E+22	6.22360E-18	4.31292E-25	2.48965E-39	6.13959E-82	7.47824-153
2.0E+22	5.60679E-18	3.94657E-25	2.30204E-39	5.70753E-82	6.97128-153
5.0E+22	4.67258E-18	3.34875E-25	1.97407E-39	4.91590E-82	6.01377-153
1.0E+23	2.96441E-18	2.16146E-25	1.28895E-39	3.23148E-82	3.96684-153
2.0E+23	1.44765E-18	1.07621E-25	6.48120E-40	1.62614E-82	1.98836-153
5.0E+23	3.74534E-19	2.86729E-26	1.74848E-40	4.38425E-83	5.32456-153
1.0E+24	8.08785E-21	6.25749E-28	3.82194E-42	9.39947E-85	1.10288-153
2.0E+24	6.83445E-22	5.20470E-29	3.13298E-43	7.54866E-86	8.60985-157
5.0E+24	6.83445E-22	5.20470E-29	3.13298E-43	7.54866E-86	8.60985-157
1.0E+25	6.83445E-22	5.20470E-29	3.13298E-43	7.54866E-86	8.60985-157
2.0E+25	6.83445E-22	5.20470E-29	3.13298E-43	7.54866E-86	8.60985-157

O2 COLUMN DENSITY

Table I(b). Total photoionization Rate Coefficients for the Process $O_3 + h\nu \rightarrow O(^1D) + O_2(^1A_g)$ for Wavelengths Below 50 nm

	O3 COLUMN DENSITY									
	1.0E+12	2.0E+12	5.0E+12	1.0E+13	2.0E+13	5.0E+13	1.0E+14	2.0E+14	5.0E+14	1.0E+15
1.0E+15	9.72657E-03	9.72631E-03	9.72597E-03	9.72530E-03	9.72328E-03	9.71992E-03	9.71321E-03	9.69311E-03	9.65972E-03	9.65972E-03
2.0E+15	9.72657E-03	9.72650E-03	9.72596E-03	9.72529E-03	9.72328E-03	9.71992E-03	9.71321E-03	9.69311E-03	9.65970E-03	9.65970E-03
5.0E+15	9.72655E-03	9.72649E-03	9.72628E-03	9.72595E-03	9.72326E-03	9.71990E-03	9.71319E-03	9.69309E-03	9.65968E-03	9.65968E-03
1.0E+16	9.72653E-03	9.72646E-03	9.72626E-03	9.72592E-03	9.72323E-03	9.71988E-03	9.71316E-03	9.69306E-03	9.65966E-03	9.65966E-03
2.0E+16	9.72648E-03	9.72641E-03	9.72587E-03	9.72520E-03	9.72319E-03	9.71983E-03	9.71312E-03	9.69302E-03	9.65953E-03	9.65953E-03
5.0E+16	9.72636E-03	9.72630E-03	9.72609E-03	9.72576E-03	9.72308E-03	9.72307E-03	9.71971E-03	9.71330E-03	9.69290E-03	9.65945E-03
1.0E+17	9.72622E-03	9.72616E-03	9.72596E-03	9.72562E-03	9.72495E-03	9.72293E-03	9.71957E-03	9.71266E-03	9.69276E-03	9.65938E-03
2.0E+17	9.72606E-03	9.72600E-03	9.72579E-03	9.72546E-03	9.72479E-03	9.72277E-03	9.71941E-03	9.71270E-03	9.69260E-03	9.65922E-03
5.0E+17	9.72585E-03	9.72579E-03	9.72559E-03	9.72526E-03	9.72458E-03	9.72257E-03	9.71921E-03	9.71250E-03	9.69240E-03	9.65902E-03
1.0E+18	9.72552E-03	9.72563E-03	9.72543E-03	9.72509E-03	9.72242E-03	9.72242E-03	9.71905E-03	9.71234E-03	9.69224E-03	9.65885E-03
2.0E+18	9.72552E-03	9.72545E-03	9.72520E-03	9.72491E-03	9.72424E-03	9.72233E-03	9.71887E-03	9.71215E-03	9.69206E-03	9.65867E-03
5.0E+18	9.72528E-03	9.72522E-03	9.72502E-03	9.72468E-03	9.72301E-03	9.72199E-03	9.71863E-03	9.71132E-03	9.69182E-03	9.65846E-03
1.0E+19	9.72514E-03	9.72507E-03	9.72487E-03	9.72453E-03	9.72386E-03	9.72184E-03	9.71849E-03	9.71178E-03	9.69168E-03	9.65830E-03
2.0E+19	9.72499E-03	9.72492E-03	9.72472E-03	9.72439E-03	9.72372E-03	9.72170E-03	9.71834E-03	9.71153E-03	9.69153E-03	9.65815E-03
5.0E+19	9.72474E-03	9.72467E-03	9.72447E-03	9.72413E-03	9.72346E-03	9.72145E-03	9.71809E-03	9.71138E-03	9.69128E-03	9.65790E-03
1.0E+20	9.72443E-03	9.72436E-03	9.72416E-03	9.72413E-03	9.72316E-03	9.72144E-03	9.71778E-03	9.71107E-03	9.69097E-03	9.65759E-03
2.0E+20	9.72394E-03	9.72387E-03	9.72367E-03	9.72333E-03	9.72266E-03	9.72065E-03	9.71729E-03	9.71058E-03	9.69048E-03	9.65710E-03
5.0E+20	9.72295E-03	9.72283E-03	9.72268E-03	9.72235E-03	9.72156E-03	9.71966E-03	9.71630E-03	9.70959E-03	9.68950E-03	9.65612E-03
1.0E+21	9.72106E-03	9.72100E-03	9.72080E-03	9.72046E-03	9.71979E-03	9.71777E-03	9.71442E-03	9.70771E-03	9.68761E-03	9.65424E-03
2.0E+21	9.71753E-03	9.71752E-03	9.71732E-03	9.71698E-03	9.71631E-03	9.71430E-03	9.71094E-03	9.70742E-03	9.68414E-03	9.65078E-03
5.0E+21	9.70753E-03	9.70752E-03	9.70736E-03	9.70703E-03	9.70636E-03	9.70434E-03	9.70099E-03	9.69429E-03	9.67421E-03	9.54087E-03
1.0E+22	9.69154E-03	9.69128E-03	9.69108E-03	9.69074E-03	9.69002E-03	9.68864E-03	9.68471E-03	9.67801E-03	9.65796E-03	9.62465E-03
2.0E+22	9.66011E-03	9.66005E-03	9.65985E-03	9.65951E-03	9.65884E-03	9.65684E-03	9.65349E-03	9.64681E-03	9.59356E-03	9.56280E-03
5.0E+22	9.57430E-03	9.57423E-03	9.57404E-03	9.57370E-03	9.57304E-03	9.57104E-03	9.56772E-03	9.56108E-03	9.54119E-03	9.50815E-03
1.0E+23	9.45231E-03	9.45224E-03	9.45205E-03	9.45172E-03	9.45106E-03	9.44908E-03	9.44579E-03	9.43921E-03	9.41949E-03	9.38676E-03
2.0E+23	9.26850E-03	9.26830E-03	9.26805E-03	9.26798E-03	9.26733E-03	9.26538E-03	9.26213E-03	9.25565E-03	9.23622E-03	9.20336E-03
5.0E+23	8.97590E-03	8.97564E-03	8.97534E-03	8.97504E-03	8.97470E-03	8.97281E-03	8.96965E-03	8.96333E-03	8.94443E-03	8.91309E-03
1.0E+24	8.79831E-03	8.79795E-03	8.79776E-03	8.79745E-03	8.79683E-03	8.79683E-03	8.79497E-03	8.78595E-03	8.76715E-03	8.73637E-03
2.0E+24	8.69314E-03	8.69308E-03	8.69290E-03	8.69259E-03	8.69198E-03	8.6915E-03	8.68709E-03	8.68095E-03	8.66271E-03	8.62235E-03
5.0E+24	8.63533E-03	8.63532E-03	8.63514E-03	8.63483E-03	8.63423E-03	8.63241E-03	8.62938E-03	8.62333E-03	8.60520E-03	8.57509E-03
1.0E+25	8.62788E-03	8.62782E-03	8.62733E-03	8.62673E-03	8.62491E-03	8.61584E-03	8.59773E-03	8.56766E-03	8.53777E-03	8.50777E-03
2.0E+25	8.62761E-03	8.62755E-03	8.62737E-03	8.62707E-03	8.62646E-03	8.62162E-03	8.61557E-03	8.61557E-03	8.59747E-03	8.56739E-03

Table B6. Total Photoassociation Rate Coefficients for the Process $O_3 + h\nu \rightarrow O(^1\Delta_g) + O_2(^1\Lambda_g)$ for Wavelengths Below 310 nm (Contd.)

	03 COLUMN DENSITY					
	2.0E+15	5.0E+15	1.0E+16	2.0E+16	5.0E+16	1.0E+17
1.0E+15	9.59335E-03	9.39780E-03	9.08291E-03	8.49245E-03	6.99530E-03	5.19826E-03
2.0E+15	9.59338E-03	9.39779E-03	9.08290E-03	8.49245E-03	6.99530E-03	5.19826E-03
3.0E+15	9.59337E-03	9.39776E-03	9.08289E-03	8.49244E-03	6.99528E-03	5.19825E-03
4.0E+15	9.59334E-03	9.39775E-03	9.08287E-03	8.49241E-03	6.99526E-03	5.19823E-03
5.0E+15	9.59324E-03	9.39770E-03	9.08282E-03	8.49237E-03	6.99523E-03	5.19821E-03
6.0E+15	9.59324E-03	9.39759E-03	9.08271E-03	8.49227E-03	6.99514E-03	5.19814E-03
7.0E+15	9.59316E-03	9.39759E-03	9.08258E-03	8.49214E-03	6.99504E-03	5.19805E-03
8.0E+15	9.59304E-03	9.39746E-03	9.08243E-03	8.49214E-03	6.99492E-03	5.19796E-03
9.0E+15	9.59298E-03	9.39730E-03	9.08243E-03	8.49200E-03	6.99492E-03	5.19796E-03
1.0E+16	9.59268E-03	9.39711E-03	9.08224E-03	8.49182E-03	6.99476E-03	5.19783E-03
2.0E+16	9.59252E-03	9.39695E-03	9.08208E-03	8.49168E-03	6.99464E-03	5.19772E-03
3.0E+16	9.59234E-03	9.39677E-03	9.08191E-03	8.49152E-03	6.99450E-03	5.19760E-03
4.0E+16	9.59211E-03	9.39655E-03	9.08170E-03	8.49133E-03	6.99432E-03	5.19745E-03
5.0E+16	9.59211E-03	9.39655E-03	9.08156E-03	8.49113E-03	6.99420E-03	5.19735E-03
6.0E+16	9.59197E-03	9.39640E-03	9.08142E-03	8.49104E-03	6.99407E-03	5.19722E-03
7.0E+16	9.59182E-03	9.39626E-03	9.08142E-03	8.49104E-03	6.99407E-03	5.19722E-03
8.0E+16	9.59157E-03	9.39601E-03	9.08131E-03	8.49080E-03	6.99384E-03	5.19702E-03
9.0E+16	9.59127E-03	9.39571E-03	9.08087E-03	8.49051E-03	6.99357E-03	5.19678E-03
1.0E+17	9.59078E-03	9.39522E-03	9.08039E-03	8.49039E-03	6.99315E-03	5.19641E-03
2.0E+17	9.59078E-03	9.39522E-03	9.08039E-03	8.49005E-03	6.99309E-03	5.19641E-03
3.0E+17	9.58983E-03	9.39525E-03	9.07945E-03	8.48915E-03	6.99236E-03	5.19677E-03
4.0E+17	9.58979E-03	9.39524E-03	9.07763E-03	8.48730E-03	6.99081E-03	5.19477E-03
5.0E+17	9.58448E-03	9.38901E-03	9.07430E-03	8.48421E-03	6.98799E-03	5.19214E-03
6.0E+17	9.57461E-03	9.37927E-03	9.06478E-03	8.47510E-03	6.97998E-03	5.18561E-03
7.0E+17	9.55847E-03	9.36335E-03	9.04921E-03	8.46020E-03	6.96689E-03	5.17494E-03
8.0E+17	9.52753E-03	9.33282E-03	9.01937E-03	8.43168E-03	6.94189E-03	5.15466E-03
9.0E+17	9.44251E-03	9.24904E-03	8.93748E-03	8.35348E-03	6.87359E-03	5.09956E-03
1.0E+18	9.32171E-03	9.12954E-03	8.82126E-03	8.24273E-03	6.77734E-03	5.02258E-03
2.0E+18	9.13988E-03	9.55889E-03	8.64678E-03	8.07623E-03	6.63473E-03	4.91022E-03
3.0E+18	8.85063E-03	9.05673E-03	8.37093E-03	7.814562E-03	6.41556E-03	4.74424E-03
4.0E+18	8.67521E-03	8.43496E-03	8.26496E-03	7.68191E-03	6.20574E-03	4.05444E-03
5.0E+18	8.57203E-03	8.36424E-03	8.16827E-03	7.57741E-03	6.19755E-03	4.60753E-03
6.0E+18	8.51528E-03	8.38932E-03	8.05539E-03	7.4433LE-03	6.18260E-03	4.58363E-03
7.0E+18	8.50791E-03	8.33232E-03	8.04850E-03	7.31806F-03	6.17782E-03	4.58058E-03
8.0E+18	8.50765E-03	8.33157E-03	8.04925E-03	7.51744E-03	6.17765E-03	4.58647E-03
9.0E+18	8.50765E-03	8.33157E-03	8.04925E-03	7.51744E-03	6.17765E-03	4.58647E-03
1.0E+19	8.50765E-03	8.33157E-03	8.04925E-03	7.51744E-03	6.17765E-03	4.58647E-03

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Table B9. Total Photodissociation rate Coefficients for the Process $O_3 + h\nu \rightarrow O(^1D) + O_2(^1\Delta_g)$ for Wavelengths Below 10 nm (Contd.)

03 COLUMN DENSITY						
	5.0E+18	1.0E+19	2.0E+19	5.0E+19	1.0E+20	2.0E+20
1.0E+15	1.09487E-04	3.98002E-05	9.65152E-06	3.31466E-07	1.71194E-09	4.71332E-14
2.0E+15	1.09487E-04	3.98002E-05	9.65152E-06	3.31466E-07	1.71194E-09	4.71332E-14
5.0E+15	1.09487E-04	3.98002E-05	9.65152E-06	3.31466E-07	1.71194E-09	4.71332E-14
1.0E+16	1.09487E-04	3.98002E-05	9.65152E-06	3.31466E-07	1.71194E-09	4.71332E-14
2.0E+16	1.09486E-04	3.98002E-05	9.65152E-06	3.31466E-07	1.71194E-09	4.71332E-14
5.0E+16	1.09486E-04	3.98002E-05	9.65152E-06	3.31466E-07	1.71194E-09	4.71332E-14
1.0E+17	1.09486E-04	3.98002E-05	9.65152E-06	3.31466E-07	1.71194E-09	4.71332E-14
2.0E+17	1.09485E-04	3.98002E-05	9.65152E-06	3.31466E-07	1.71194E-09	4.71332E-14
5.0E+17	1.09485E-04	3.98002E-05	9.65152E-06	3.31466E-07	1.71194E-09	4.71332E-14
1.0E+18	1.09482E-04	3.98002E-05	9.65152E-06	3.31466E-07	1.71194E-09	4.71332E-14
2.0E+18	1.09480E-04	3.98002E-05	9.65152E-06	3.31466E-07	1.71194E-09	4.71332E-14
5.0E+18	1.09477E-04	3.98002E-05	9.65152E-06	3.31466E-07	1.71194E-09	4.71332E-14
1.0E+19	1.09475E-04	3.97444E-05	9.65152E-06	3.31466E-07	1.71194E-09	4.71332E-14
2.0E+19	1.09472E-04	3.97343E-05	9.65152E-06	3.31466E-07	1.71194E-09	4.71332E-14
5.0E+19	1.09456E-04	3.97934E-05	9.65151E-06	3.31466E-07	1.71194E-09	4.71332E-14
1.0E+20	1.09450E-04	3.97889E-05	9.65150E-06	3.31466E-07	1.71194E-09	4.71332E-14
2.0E+20	1.09449E-04	3.97982E-05	9.65149E-06	3.31466E-07	1.71194E-09	4.71332E-14
5.0E+20	1.09449E-04	3.97980E-05	9.65149E-06	3.31466E-07	1.71194E-09	4.71332E-14
1.0E+21	1.09447E-04	3.97961E-05	9.65146E-06	3.31466E-07	1.71194E-09	4.71332E-14
2.0E+21	1.09447E-04	3.97935E-05	9.65141E-06	3.31466E-07	1.71194E-09	4.71332E-14
5.0E+21	1.09447E-04	3.97885E-05	9.65131E-06	3.31466E-07	1.71194E-09	4.71332E-14
1.0E+22	1.09447E-04	3.97792E-05	9.65111E-06	3.31466E-07	1.71194E-09	4.71332E-14
2.0E+22	1.09447E-04	3.97652E-05	9.65080E-06	3.31466E-07	1.71194E-09	4.71332E-14
5.0E+22	1.09447E-04	3.97372E-05	9.65022E-06	3.31466E-07	1.71194E-09	4.71332E-14
1.0E+23	1.08418E-04	3.96965E-05	9.64923E-06	3.31466E-07	1.71194E-09	4.71332E-14
2.0E+23	1.08417E-04	3.96586E-05	9.64832E-06	3.31466E-07	1.71194E-09	4.71332E-14
5.0E+23	1.07911E-04	3.96340E-05	9.64769E-06	3.31466E-07	1.71194E-09	4.71332E-14
1.0E+24	1.07845E-04	3.96286E-05	9.64750E-06	3.31466E-07	1.71194E-09	4.71332E-14
2.0E+24	1.07844E-04	3.96285E-05	9.64750E-06	3.31466E-07	1.71194E-09	4.71332E-14
5.0E+24	1.07844E-04	3.96285E-05	9.64750E-06	3.31466E-07	1.71194E-09	4.71332E-14
1.0E+25	1.07844E-04	3.96285E-05	9.64750E-06	3.31466E-07	1.71194E-09	4.71332E-14
2.0E+25	1.07844E-04	3.96285E-05	9.64750E-06	3.31466E-07	1.71194E-09	4.71332E-14

Table B7. Total Photodissociation Rate Coefficients for the Process $O_3 + h\nu \rightarrow O + O_2$ for Wavelengths From 310 to 730 nm

	O3 COLUMN DENSITY
1.0E+14	4.17996E-04
2.0E+14	4.17995E-04
5.0E+14	4.17994E-04
1.0E+15	4.17992E-04
2.0E+15	4.17989E-04
5.0E+15	4.17978E-04
1.0E+16	4.17960E-04
2.0E+16	4.17924E-04
5.0E+16	4.17816E-04
1.0E+17	4.17636E-04
2.0E+17	4.17276E-04
5.0E+17	4.16205E-04
1.0E+18	4.14441E-04
2.0E+18	4.10992E-04
5.0E+18	4.01237E-04
1.0E+19	3.86702E-04
2.0E+19	3.62634E-04
5.0E+19	3.13047E-04
1.0E+20	2.60086E-04
2.0E+20	1.90049E-04
5.0E+20	8.70967E-05
1.0E+21	3.45505E-05

Table B6. Total Photodissociation Rate Coefficients for the Process $H_2O + h\nu \rightarrow OH + H$ for Wavelengths From 121.9 to 186.4 nm

03 COLUMN DENSITY													
	1.0E+11	2.0E+11	5.0E+11	1.0E+12	2.0E+12	5.0E+12	1.0E+13	2.0E+13	5.0E+13	1.0E+14	2.0E+14	5.0E+14	
1.0E+13	7.38657E-06	7.38656E-06	7.38653E-06	7.38647E-06	7.38637E-06	7.38606E-06	7.38554E-06	7.38451E-06	7.38141E-06	7.37625E-06	7.37425E-06	7.37245E-06	
2.0E+13	7.38653E-06	7.38652E-06	7.38649E-06	7.38643E-06	7.38602E-06	7.38550E-06	7.38447E-06	7.38137E-06	7.37621E-06	7.37421E-06	7.37221E-06	7.37021E-06	
5.0E+13	7.38640E-06	7.38639E-06	7.38636E-06	7.38631E-06	7.38621E-06	7.38590E-06	7.38435E-06	7.38125E-06	7.37609E-06	7.37409E-06	7.37209E-06	7.37009E-06	
1.0E+14	7.38620E-06	7.38619E-06	7.38616E-06	7.38611E-06	7.38600E-06	7.38569E-06	7.38518E-06	7.38414E-06	7.38104E-06	7.37588E-06	7.37488E-06	7.37288E-06	
2.0E+14	7.38579E-06	7.38578E-06	7.38575E-06	7.38570E-06	7.38560E-06	7.38529E-06	7.38477E-06	7.38374E-06	7.38064E-06	7.37547E-06	7.37447E-06	7.37247E-06	
5.0E+14	7.38457E-06	7.38456E-06	7.38453E-06	7.38448E-06	7.38443E-06	7.38406E-06	7.38405E-06	7.38251E-06	7.37941E-06	7.37425E-06	7.37225E-06	7.37025E-06	
1.0E+15	7.38254E-06	7.38253E-06	7.38250E-06	7.38245E-06	7.38234E-06	7.38203E-06	7.38152E-06	7.38048E-06	7.37738E-06	7.37222E-06	7.37022E-06	7.36822E-06	
2.0E+15	7.37849E-06	7.37848E-06	7.37845E-06	7.37840E-06	7.37834E-06	7.37798E-06	7.37747E-06	7.37643E-06	7.37333E-06	7.36817E-06	7.36617E-06	7.36417E-06	
5.0E+15	7.36647E-06	7.36646E-06	7.36643E-06	7.36638E-06	7.36633E-06	7.36596E-06	7.36545E-06	7.36441E-06	7.36132E-06	7.35616E-06	7.35316E-06	7.35016E-06	
1.0E+16	7.34685E-06	7.34684E-06	7.34683E-06	7.34682E-06	7.34665E-06	7.34663E-06	7.34652E-06	7.34479E-06	7.34169E-06	7.33654E-06	7.33254E-06	7.32854E-06	
2.0E+16	7.30904E-06	7.30903E-06	7.30895E-06	7.30885E-06	7.30885E-06	7.30854E-06	7.30854E-06	7.30390E-06	7.30390E-06	7.29875E-06	7.29875E-06	7.29477E-06	
5.0E+16	7.20579E-06	7.20577E-06	7.20574E-06	7.20574E-06	7.20559E-06	7.20559E-06	7.20528E-06	7.20374E-06	7.20065E-06	7.19552E-06	7.19552E-06	7.19152E-06	
1.0E+17	7.06019E-06	7.06018E-06	7.06015E-06	7.06015E-06	7.06010E-06	7.06009E-06	7.05996E-06	7.05918E-06	7.05508E-06	7.04996E-06	7.04996E-06	7.04596E-06	
2.0E+17	6.82192E-06	6.82191E-06	6.82188E-06	6.82183E-06	6.82172E-06	6.82142E-06	6.82091E-06	6.81989E-06	6.81683E-06	6.81174E-06	6.80774E-06	6.80374E-06	
5.0E+17	6.35730E-06	6.35729E-06	6.35726E-06	6.35721E-06	6.35711E-06	6.35681E-06	6.35631E-06	6.35530E-06	6.35522E-06	6.34723E-06	6.34223E-06	6.34223E-06	
1.0E+18	5.89929E-06	5.89928E-06	5.89927E-06	5.89926E-06	5.89910E-06	5.89880E-06	5.89830E-06	5.89730E-06	5.89431E-06	5.88934E-06	5.88934E-06	5.88934E-06	
2.0E+18	5.40915E-06	5.40914E-06	5.40912E-06	5.40912E-06	5.40907E-06	5.40907E-06	5.40867E-06	5.40819E-06	5.40427E-06	5.39939E-06	5.39939E-06	5.39939E-06	
5.0E+18	4.82323E-06	4.82322E-06	4.82319E-06	4.82316E-06	4.82305E-06	4.82277E-06	4.82230E-06	4.82136E-06	4.81854E-06	4.81395E-06	4.81395E-06	4.81395E-06	
1.0E+19	4.41601E-06	4.41600E-06	4.41598E-06	4.41593E-06	4.41584E-06	4.41558E-06	4.41513E-06	4.41424E-06	4.41158E-06	4.40715E-06	4.40315E-06	4.40315E-06	
2.0E+19	3.91448E-06	3.91447E-06	3.91445E-06	3.91443E-06	3.91433E-06	3.91409E-06	3.91369E-06	3.91289E-06	3.91048E-06	3.90648E-06	3.90648E-06	3.90648E-06	
5.0E+19	2.88069E-06	2.88069E-06	2.88067E-06	2.88064E-06	2.88058E-06	2.88052E-06	2.88040E-06	2.88011E-06	2.87951E-06	2.87773E-06	2.87477E-06	2.87477E-06	
1.0E+20	1.78257E-06	1.78256E-06	1.78254E-06	1.78252E-06	1.78250E-06	1.78239E-06	1.78221E-06	1.78185E-06	1.78077E-06	1.77897E-06	1.77897E-06	1.77897E-06	
2.0E+20	7.20230E-07	7.20229E-07	7.20224E-07	7.20218E-07	7.20204E-07	7.20165E-07	7.20098E-07	7.19965E-07	7.19566E-07	7.18902E-07	7.18902E-07	7.18902E-07	
5.0E+20	1.80122E-07	1.80122E-07	1.80121E-07	1.80121E-07	1.80121E-07	1.80118E-07	1.80114E-07	1.80107E-07	1.80084E-07	1.80045E-07	1.80045E-07	1.80045E-07	
1.0E+21	6.60228E-08	6.60228E-08	6.60227E-08	6.60227E-08	6.60227E-08	6.60227E-08	6.60225E-08	6.60222E-08	6.60217E-08	6.60174E-08	6.60174E-08	6.60174E-08	
2.0E+21	2.82041E-08	2.82041E-08	2.82040E-08	2.82040E-08	2.82040E-08	2.82039E-08	2.82036E-08	2.82030E-08	2.82020E-08	2.82020E-08	2.82020E-08	2.82020E-08	
5.0E+21	1.06830E-08	1.06829E-08	1.06823E-08	1.06823E-08	1.06823E-08	1.06823E-08							
1.0E+22	2.28878E-09	2.28878E-09	2.28878E-09	2.28878E-09	2.28878E-09	2.28878E-09	2.28877E-09	2.28875E-09	2.28862E-09	2.28862E-09	2.28862E-09	2.28862E-09	
2.0E+22	6.07985E-10	6.07985E-10	6.07985E-10	6.07985E-10	6.07985E-10	6.07985E-10	6.07984E-10	6.07983E-10	6.07983E-10	6.07940E-10	6.07940E-10	6.07940E-10	
5.0E+22	3.05080E-10	3.05079E-10	3.05079E-10	3.05056E-10	3.05056E-10	3.05056E-10							
1.0E+23	2.82473E-10	2.82473E-10	2.82473E-10	2.82473E-10	2.82473E-10	2.82473E-10	2.82472E-10	2.82471E-10	2.82462E-10	2.82462E-10	2.82462E-10	2.82462E-10	

Tab. 133. Total Photodissociation $\text{H}_2\text{O} + \text{h}\nu \rightarrow \text{OH} + \text{H}$ for Wavelengths From 121.9 to 166.4 nm (Cont'd)

03 COLUMN DENSITY						
	2.0E+14	5.0E+14	1.0E+15	2.0E+15	5.0E+15	1.0E+16
1.0E+13	7.36594E-06	7.33516E-06	7.18433E-06	7.18434E-06	6.89750E-06	6.46021E-06
2.0E+13	7.36590E-06	7.33512E-06	7.18421E-06	7.18430E-06	6.89746E-06	6.46017E-06
3.0E+13	7.36578E-06	7.33500CE-06	7.1841bE-06	7.1841BE-06	6.89734E-06	6.46005E-06
4.0E+13	7.36556E-06	7.33490CE-06	7.1841bE-06	7.1841BE-06	6.89734E-06	6.46005E-06
5.0E+13	7.36555E-06	7.33480E-06	7.18396E-06	7.18398E-06	6.89714E-06	6.45985E-06
6.0E+13	7.36555E-06	7.33439E-06	7.18357E-06	7.18357E-06	6.89673E-06	6.45946E-06
7.0E+13	7.36555E-06	7.33439E-06	7.18233E-06	7.18236E-06	6.89553E-06	6.45827E-06
8.0E+13	7.36312E-06	7.33317E-06	7.18034E-06	7.18034E-06	6.89533E-06	6.45629E-06
9.0E+13	7.36142E-06	7.33114E-06	7.17631E-06	7.17631E-06	6.89533E-06	6.45564E-06
1.0E+14	7.35787E-06	7.32710E-06	7.27627E-06	7.27627E-06	6.89533E-06	6.45524E-06
2.0E+15	7.34566E-06	7.31510E-06	7.26429E-06	7.16436E-06	6.87768E-06	6.44066E-06
3.0E+15	7.32624E-06	7.29550E-06	7.24472E-06	7.14484E-06	6.85833E-06	6.42157E-06
4.0E+15	7.28846E-06	7.25775E-06	7.20702E-06	7.10725E-06	6.82104E-06	6.38477E-06
5.0E+15	7.18526E-06	7.15463E-06	7.104C3E-06	7.00454E-06	6.71913E-06	6.28416E-06
6.0E+15	7.03974E-06	7.00922E-06	6.35880E-06	6.85967E-06	6.57532E-06	6.14202E-06
7.0E+15	6.80157E-06	6.77121E-06	6.72107E-06	6.62247E-06	6.33970E-06	5.90896E-06
8.0E+15	6.33717E-06	6.30713E-06	6.25752E-06	6.15996E-06	5.88026E-06	5.45437E-06
9.0E+15	5.87940E-05	5.84973E-06	5.80072E-06	5.70437E-06	5.42617E-06	5.0780E-06
1.0E+16	5.38963E-06	5.36050E-06	5.31239E-06	5.21782E-06	4.94676E-06	4.53438E-06
2.0E+16	4.80449E-06	4.77654E-06	4.73038E-06	4.63963E-06	4.37960E-06	4.08449E-06
3.0E+16	4.39830E-06	4.37187E-06	4.32822E-06	4.24242E-06	3.90896E-06	3.62226E-06
4.0E+16	3.89849E-06	3.87464E-06	3.83524E-06	3.75781E-06	3.53555E-06	3.19886E-06
5.0E+16	2.86888E-06	2.85119E-06	2.82202E-06	2.76469E-06	2.60042E-06	2.35070E-06
6.0E+16	1.77530E-06	1.76466E-06	1.74696E-06	1.71216E-06	1.61244E-06	1.46066E-06
7.0E+16	7.17577E-07	7.13618E-07	7.07081E-07	6.94231E-07	6.57405E-07	5.06945E-07
8.0E+16	1.79968E-07	1.79379E-07	1.79359E-07	1.78610E-07	1.76451E-07	1.67344E-07
9.0E+16	6.60121E-08	6.59962E-08	6.59696E-08	6.59566E-08	6.57585E-08	6.54974E-08
1.0E+17	2.82000E-08	2.81936E-08	2.81836E-08	2.81632E-08	2.81021E-08	2.80004E-08
2.0E+17	1.06815E-09	1.06793E-08	1.06755E-08	1.06579E-08	1.06452E-08	1.06075E-08
3.0E+17	2.28845E-03	2.28797E-09	2.28717E-09	2.28071E-09	2.27267E-09	2.25667E-09
4.0E+17	6.07896E-10	6.07761E-10	6.07537E-10	6.07989E-10	6.05748E-10	5.99098E-10
5.0E+17	3.05022E-10	3.04900E-10	3.04840E-10	3.04599E-10	3.02684E-10	3.00306E-10
6.0E+17	2.82429E-10	2.82360E-10	2.82248E-10	2.82023E-10	2.81349E-10	2.78003E-10
7.0E+17	2.82429E-10	2.82360E-10	2.82248E-10	2.82023E-10	2.81349E-10	2.78003E-10

Table 186. Total Photoionisation Rate Coefficients for the Process $\text{H}_2\text{O} + \text{h}\nu \rightarrow \text{OH} + \text{H}$ for Wavelengths From 121.9 to 186.4 nm (Cont'd)

	03 COLUMN DENSITY					
	5.0E+17	1.0E+18	2.0E+18	5.0E+18	1.0E+19	2.0E+19
1.0E+13	1.92246E-06	1.25722E-06	5.46968E-07	4.69360E-08	8.30989E-10	3.13497E-13
2.0E+13	1.92245E-06	1.25721E-06	5.46964E-07	4.69357E-08	8.30985E-10	3.13496E-13
5.0E+13	1.92239E-06	1.25717E-06	5.46952E-07	4.69349E-08	8.30972E-10	3.13493E-13
1.0E+14	1.92230E-06	1.25712E-06	5.46941E-07	4.69334E-08	8.30952E-10	3.13488E-13
2.0E+14	1.92211E-06	1.25701E-06	5.46899CE-07	4.69305E-08	8.30911E-10	3.13478E-13
5.0E+14	1.92156E-06	1.25669E-06	5.46767E-07	4.69219E-08	8.30788E-10	3.13449E-13
1.0E+15	1.92063E-06	1.25615E-06	5.46561E-07	4.69075E-08	8.30584E-10	3.13400E-13
2.0E+15	1.91878E-C6	1.25507E-06	5.46151E-07	4.68787E-08	8.30176E-10	3.13303E-13
5.0E+15	1.91328E-06	1.25186E-06	5.43326E-07	4.67928E-08	8.28556E-10	3.13012E-13
1.0E+16	1.90422E-06	1.24657E-06	5.42901E-07	4.66504E-08	8.26930E-10	3.12529E-13
2.0E+16	1.88754E-06	1.23619E-06	5.38514E-07	4.63691E-08	8.22923E-10	3.11572E-13
5.0E+16	1.83662E-06	1.20658E-06	5.27438E-07	4.55517E-08	8.11219E-10	3.08767E-13
1.0E+17	1.76241E-06	1.16175E-06	5.09773E-07	4.42723E-08	7.92723E-10	3.04302E-13
2.0E+17	1.63141E-06	1.08022E-06	4.76607E-07	4.17655E-08	7.55026E-10	2.94565E-13
5.0E+17	1.36039E-06	9.06076E-07	4.03288E-07	3.59825E-08	6.65582E-10	2.70930E-13
1.0E+18	1.09022E-06	7.28659E-07	3.26428E-07	2.96381E-08	5.63528E-10	2.42722E-13
2.0E+18	8.14274E-07	5.46233E-07	2.46519E-07	2.28901E-08	4.52332E-10	2.10869E-13
5.0E+18	5.36387E-07	3.62213E-07	1.65588E-07	1.59748E-08	3.55915E-10	1.75261E-13
1.0E+19	4.10798E-07	2.78686E-07	1.28502E-07	1.27069E-08	2.77544E-10	1.54356E-13
2.0E+19	3.22841E-07	2.19649E-07	1.01880E-07	1.02443E-08	2.40483E-09	1.34277E-13
5.0E+19	2.27430E-07	1.55232E-07	7.24053E-07	7.41035E-09	1.71387E-10	1.06409E-13
1.0E+20	1.62437E-07	1.11127E-07	5.20745E-08	5.40725E-09	1.28313E-10	8.40530E-14
2.0E+20	1.04156E-07	7.14751E-08	3.37009E-08	3.56743E-09	8.75905E-11	6.15006E-14
5.0E+20	1.04151E-07	7.14751E-08	3.37009E-08	3.56743E-09	8.75905E-11	6.15006E-14
1.0E+21	4.54621E-08	3.14083E-08	1.50104E-08	1.65530E-09	4.35838E-11	3.47797E-14
2.0E+21	1.96353E-08	1.36752E-08	6.64109E-08	6.67802E-09	2.17352E-11	1.94766E-14
5.0E+21	7.49321E-09	5.25744E-09	2.59C52E-09	3.12148E-10	9.39104E-12	9.21005E-15
1.0E+22	1.60809E-09	1.13033E-09	5.59159E-10	6.8329E-11	2.11332E-12	2.19137E-15
2.0E+22	4.20746E-10	2.91428E-10	1.40184E-10	1.59283E-11	4.52356E-13	4.37401E-16
5.0E+22	2.05689E-10	1.38726E-10	6.31726E-11	6.02498E-12	1.25114E-13	6.79248E-17
1.0E+23	1.89593E-10	1.27265E-10	5.73596E-11	5.26449E-12	9.92411E-14	3.69711E-17

02 COLUMN DENSITY

Table 13. Total Ionization Rate Coefficients for the Process $\text{H}_2\text{O}_2 + \text{h}\nu \rightarrow \text{OH} + \text{OH}$ for Wavelengths From 188.2 to 303 nm

03 COLUMN DENSITY						
	1.0E+12	2.0E+12	5.0E+12	1.0E+13	2.0E+13	5.0E+13
1.0E+17	1.59387E-04	1.59386E-04	1.59384E-04	1.59380E-04	1.59373E-04	1.59351E-04
2.0E+17	1.59386E-04	1.59386E-04	1.59384E-04	1.59380E-04	1.59373E-04	1.59350E-04
5.0E+17	1.59386E-04	1.59386E-04	1.59383E-04	1.59379E-04	1.59372E-04	1.59335E-04
1.0E+18	1.59385E-04	1.59384E-04	1.59382E-04	1.59378E-04	1.59371E-04	1.59349E-04
2.0E+18	1.59382E-04	1.59382E-04	1.59380E-04	1.59376E-04	1.59368E-04	1.59346E-04
5.0E+18	1.59376E-04	1.59373E-04	1.59369E-04	1.59367E-04	1.59362E-04	1.59310E-04
1.0E+19	1.59366E-04	1.59365E-04	1.59363E-04	1.59362E-04	1.59362E-04	1.59340E-04
2.0E+19	1.59347E-04	1.59346E-04	1.59344E-04	1.59340E-04	1.59332E-04	1.59311E-04
5.0E+19	1.59236E-04	1.59230E-04	1.59229E-04	1.59229E-04	1.59226E-04	1.59225E-04
1.0E+20	1.59230E-04	1.59229E-04	1.59227E-04	1.59223E-04	1.59216E-04	1.59194E-04
2.0E+20	1.59117E-04	1.59116E-04	1.59114E-04	1.59110E-04	1.59103E-04	1.59081E-04
5.0E+20	1.59015E-04	1.59012E-04	1.59010E-04	1.59007E-04	1.59001E-04	1.58971E-04
1.0E+21	1.58681E-04	1.58680E-04	1.58678E-04	1.58674E-04	1.58667E-04	1.58660E-04
2.0E+21	1.58161E-04	1.58160E-04	1.58158E-04	1.58154E-04	1.58147E-04	1.58125E-04
5.0E+21	1.56917E-04	1.56916E-04	1.56914E-04	1.56911E-04	1.56903E-04	1.56881E-04
1.0E+22	1.54846E-04	1.54845E-04	1.54843E-04	1.54839E-04	1.54832E-04	1.54810E-04
2.0E+22	1.51283E-04	1.51282E-04	1.51280E-04	1.51277E-04	1.51269E-04	1.51248E-04
5.0E+22	1.42514E-04	1.42513E-04	1.42511E-04	1.42508E-04	1.42501E-04	1.42480E-04
1.0E+23	1.30898E-04	1.30897E-04	1.30895E-04	1.30891E-04	1.30885E-04	1.30865E-04
2.0E+23	1.16527E-04	1.16527E-04	1.16525E-04	1.16522E-04	1.16515E-04	1.16496E-04
5.0E+23	1.00617E-04	1.00616E-04	1.00614E-04	1.00612E-04	1.00606E-04	1.00589E-04
1.0E+24	9.47553E-05	9.47547E-05	9.47531E-05	9.47504E-05	9.47450E-05	9.47288E-05
2.0E+24	9.25133E-05	9.25132E-05	9.25117E-05	9.25091E-05	9.25038E-05	9.24882E-05
5.0E+24	9.15194E-05	9.15188E-05	9.15173E-05	9.15147E-05	9.15096E-05	9.14942E-05
1.0E+25	9.13956E-05	9.13951E-05	9.13936E-05	9.13910E-05	9.13859E-05	9.13449E-05
2.0E+25	9.13912E-05	9.13907E-05	9.13892E-05	9.13866E-05	9.13815E-05	9.13661E-05
5.0E+25	9.13912E-05	9.13907E-05	9.13892E-05	9.13866E-05	9.13815E-05	9.13661E-05

02 COLUMN DENSITY

02 COLUMN DENSITY						
	1.0E+14	2.0E+14	5.0E+14	1.0E+15	2.0E+15	5.0E+15
1.0E+17	1.59314E-04	1.59314E-04	1.59314E-04	1.59314E-04	1.59314E-04	1.59314E-04
2.0E+17	1.59240E-04	1.59240E-04	1.59240E-04	1.59240E-04	1.59240E-04	1.59240E-04
5.0E+17	1.58655E-04	1.58655E-04	1.58655E-04	1.58655E-04	1.58655E-04	1.58655E-04
1.0E+18	1.58654E-04	1.58654E-04	1.58654E-04	1.58654E-04	1.58654E-04	1.58654E-04
2.0E+18	1.58633E-04	1.58633E-04	1.58633E-04	1.58633E-04	1.58633E-04	1.58633E-04
5.0E+18	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04
1.0E+19	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04
2.0E+19	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04
5.0E+19	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04
1.0E+20	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04
2.0E+20	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04
5.0E+20	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04
1.0E+21	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04
2.0E+21	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04
5.0E+21	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04
1.0E+22	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04
2.0E+22	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04
5.0E+22	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04
1.0E+23	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04
2.0E+23	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04
5.0E+23	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04
1.0E+24	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04
2.0E+24	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04
5.0E+24	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04
1.0E+25	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04
2.0E+25	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04
5.0E+25	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04	1.58632E-04

01 COLUMN DENSITY

Table B. Total Photoionisation Rate Coefficients for the Process $H_2O_2 + h\nu \rightarrow OH + OH$ for Wavelengths From 188.2 to 310 nm (Cont'd)

03 COLUMN DENSITY							
2.0E+15	5.0E+15	1.0E+16	2.0E+16	5.0E+17	2.0E+17	5.0E+18	2.0E+18
1.0E+17	1.557829E-04	1.557829E-04	1.52310E-04	1.45746E-04	1.28723E-04	1.07248E-04	8.02864E-05
2.0E+17	1.57928E-04	1.55781E-04	1.52210E-04	1.45746E-04	1.28723E-04	1.07248E-04	8.02862E-05
5.0E+17	1.57928E-04	1.55781E-04	1.52109E-04	1.45746E-04	1.28722E-04	1.07247E-04	8.02856E-05
1.0E+18	1.57927E-04	1.55779E-04	1.52308E-04	1.45744E-04	1.28721E-04	1.07246E-04	8.02846E-05
2.0E+18	1.57924E-04	1.55777E-04	1.52306E-04	1.45742E-04	1.28719E-04	1.07247E-04	8.02827E-05
5.0E+18	1.57918E-04	1.55771E-04	1.52299E-04	1.45736E-04	1.28713E-04	1.07238E-04	8.02770E-05
1.0E+19	1.57908E-04	1.55761E-04	1.52289E-04	1.45726E-04	1.28703E-04	1.07228E-04	8.02681E-05
2.0E+19	1.57830E-04	1.55742E-04	1.52270E-04	1.45707E-04	1.28685E-04	1.07211E-04	8.02516E-05
5.0E+19	1.57840E-04	1.55693E-04	1.52222E-04	1.45659E-04	1.28638E-04	1.07166E-04	8.02095E-05
1.0E+20	1.55625E-04	1.52154E-04	1.45592E-04	1.43358E-04	1.28573E-04	1.07103E-04	8.01518E-05
2.0E+20	1.57559E-04	1.55513E-04	1.52043E-04	1.45482E-04	1.28465E-04	1.07001E-04	8.00571E-05
5.0E+20	1.57558E-04	1.55412E-04	1.51943E-04	1.45384E-04	1.28374E-04	1.06918E-04	7.99879E-05
1.0E+21	1.57225E-04	1.55081E-04	1.51614E-04	1.45059E-04	1.28061E-04	1.06623E-04	7.97232E-05
2.0E+21	1.56705E-04	1.54565E-04	1.51102E-04	1.44555E-04	1.27560E-04	1.06175E-04	7.93277E-05
5.0E+21	1.55467E-04	1.53333E-04	1.49882E-04	1.43358E-04	1.26445E-04	1.05132E-04	7.84280E-05
1.0E+22	1.53404E-04	1.51281E-04	1.47849E-04	1.41362E-04	1.24553E-04	1.04211E-04	7.69211E-05
2.0E+22	1.49856E-04	1.47755E-04	1.44359E-04	1.37942E-04	1.21327E-04	1.00443E-04	7.44119E-05
5.0E+22	1.41112E-04	1.39085E-04	1.35785E-04	1.29557E-04	1.13463E-04	9.33263E-05	6.84478E-05
1.0E+23	1.29566E-04	1.27606E-04	1.24442E-04	1.18474E-04	1.03104E-04	8.39956E-05	6.06861E-05
2.0E+23	1.15275E-04	1.13435E-04	1.10466E-04	1.04876E-04	9.05393E-05	7.28821E-05	5.17791E-05
5.0E+23	9.94636E-05	9.73186E-05	9.51356E-05	9.00952E-05	7.72450E-05	6.16097E-05	4.32964E-05
1.0E+24	9.36864E-05	9.21162E-05	8.95869E-05	8.48393E-05	7.27602E-05	5.81211E-05	4.10755E-05
2.0E+24	9.14786E-05	8.910582E-05	8.75092E-05	8.29128E-05	7.12219E-05	5.70570E-05	4.05524E-05
5.0E+24	9.05015E-05	8.91064E-05	8.65982E-05	8.20783E-05	7.15801E-05	5.66424E-05	4.03767E-05
1.0E+25	9.03800E-05	8.81881E-05	8.64851E-05	8.19749E-05	7.05012E-05	5.65921E-05	4.03582E-05
2.0E+25	9.03777E-05	8.80839E-05	8.64811E-05	8.19713E-05	7.04984E-05	5.65903E-05	4.03575E-05
5.0E+25	9.03757E-05	8.80839E-05	8.64811E-05	8.19713E-05	7.04984E-05	5.65903E-05	4.03575E-05

Table 1b. Total Photo-dissociation Rate Coefficients for the Process $\text{H}_2\text{O}_2 + h\nu \rightarrow \text{OH} + \text{OH}$ for Wavelet 1, from 1900, 2 to 3000 (Continued)

	O2 COLUMN DENSITY					
	5.0E+18	1.0E+19	2.0E+19	5.0E+19	1.0E+20	2.0E+20
1.0E+17	3.06840E-06	3.96475E-07	1.00025E-08	3.00632E-08	1.723372E-13	9.76796E-20
2.0E+17	3.06839E-06	3.96494E-07	1.00025E-08	3.00632E-08	1.723372E-13	9.76795E-20
5.0E+17	3.06834E-06	3.98469E-07	1.00024E-08	3.00630E-08	1.723371E-13	9.76790E-20
1.0E+18	3.06825E-06	3.98480E-07	1.00023E-08	3.00628E-08	1.72369E-13	9.76782E-20
2.0E+18	3.05808E-06	3.98464E-07	1.00021E-08	3.00623E-08	1.72366E-13	9.76765E-20
5.0E+18	3.06758E-06	3.98416E-07	1.00014E-08	3.00607E-08	1.72357E-13	9.76717E-20
1.0E+19	3.06679E-06	3.98339E-07	1.00003E-08	3.00582E-08	1.72341E-13	9.76635E-20
2.0E+19	3.06532E-06	3.98192E-07	9.99808E-09	3.00531E-09	1.72311E-13	9.76473E-20
5.0E+19	3.06150E-06	3.97797E-07	9.99183E-09	3.00381E-09	1.72219E-13	9.75986E-20
1.0E+20	3.05619E-06	3.97223E-07	9.98207E-09	3.00134E-09	1.72067E-13	9.75178E-20
2.0E+20	3.04744E-06	3.96238E-07	9.96411E-09	2.99651E-09	1.71767E-13	9.73577E-20
5.0E+20	3.04502E-06	3.95999E-07	9.96032E-09	2.99599E-09	1.71756E-13	9.73571E-20
1.0E+21	3.04346E-06	3.93483E-07	9.91092E-09	2.98192E-09	1.70811E-13	9.68854E-20
2.0E+21	2.99480E-06	3.90040E-07	9.839867E-09	2.95988E-09	1.69486E-13	9.61232E-20
5.0E+21	2.94191E-06	3.83712E-07	9.70263E-09	2.91766E-09	1.66824E-13	9.46598E-20
1.0E+22	2.84524E-06	3.71293E-07	9.40340E-09	2.81018E-09	1.59618E-13	9.05682E-20
2.0E+22	2.70660E-06	3.53507E-07	8.96473E-09	2.64835E-09	1.48690E-13	8.43140E-20
5.0E+22	2.43341E-06	3.19813E-07	8.15846E-09	2.35940E-09	1.29412E-13	7.32892E-20
1.0E+23	2.04965E-06	2.67769E-07	6.73786E-09	1.76935E-09	8.7625E-13	4.87091E-21
2.0E+23	1.69456E-06	2.20257E-07	5.43161E-09	1.21507E-09	4.78734E-13	2.51090E-21
5.0E+23	1.46600E-06	1.88854E-07	4.52037E-09	8.03634E-14	1.75909E-14	6.90188E-21
1.0E+24	1.42056E-06	1.80997E-07	4.24356E-09	6.66642E-14	6.63379E-22	2.12690E-22
2.0E+24	1.41973E-06	1.80846E-07	4.23803E-09	6.57690E-14	6.40503E-22	2.27530E-22
5.0E+24	1.41974E-06	1.80846E-07	4.23803E-09	6.57690E-14	6.40503E-22	2.27530E-22
1.0E+25	1.41973E-06	1.80846E-07	4.23803E-09	6.57690E-14	6.40503E-22	2.27530E-22
2.0E+25	1.41973E-06	1.80846E-07	4.23803E-09	6.57690E-14	6.40503E-22	2.27530E-22
5.0E+25	1.41973E-06	1.80846E-07	4.23803E-09	6.57690E-14	6.40503E-22	2.27530E-22

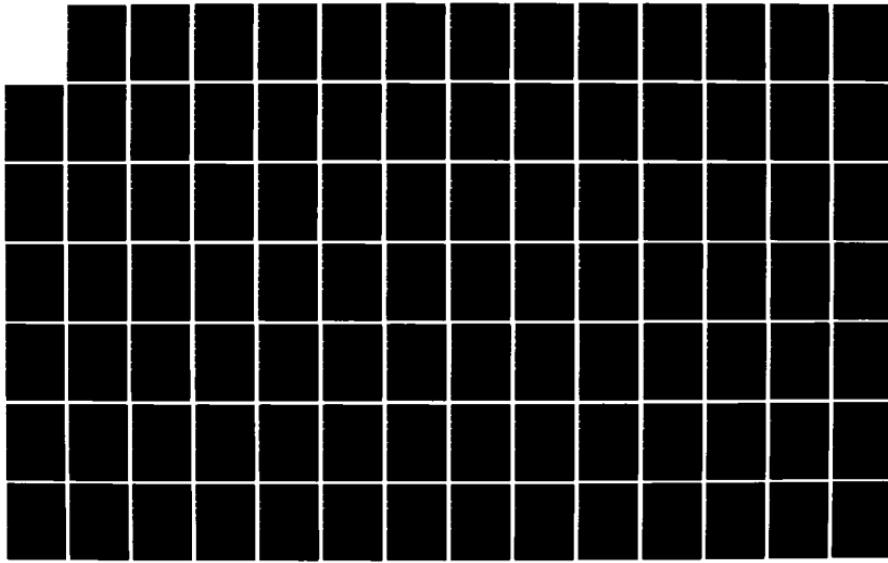
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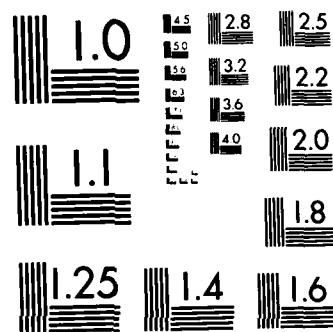
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Table B10. Total Photodissociation Rate Coefficients for the Process $\text{NO}_2 + h\nu \rightarrow \text{NO} + \text{O}^{\text{(1D)}}$ for Wavelengths From 190 to 420 nm

	1.0E+14	2.0E+14	5.0E+14	1.0E+15	2.0E+15	5.0E+15	1.0E+16	2.0E+16	5.0E+16	1.0E+17
	03 COLUMN DENSITY									
	02 COLUMN DENSITY									
1.0E+19	1.51038E-02	1.51038E-02	1.51036E-02	1.51033E-02	1.51027E-02	1.51008E-02	1.50920E-02	1.50759E-02	1.50530E-02	1.50530E-02
2.0E+19	1.51038E-02	1.51038E-02	1.51036E-02	1.51033E-02	1.51026E-02	1.51008E-02	1.50920E-02	1.50759E-02	1.50530E-02	1.50530E-02
5.0E+19	1.51038E-02	1.51037E-02	1.51035E-02	1.51032E-02	1.51026E-02	1.51008E-02	1.50920E-02	1.50759E-02	1.50529E-02	1.50529E-02
1.0E+20	1.51037E-02	1.51037E-02	1.51035E-02	1.51032E-02	1.51026E-02	1.51007E-02	1.50919E-02	1.50758E-02	1.50529E-02	1.50529E-02
2.0E+20	1.51037E-02	1.51036E-02	1.51034E-02	1.51031E-02	1.51025E-02	1.51007E-02	1.50918E-02	1.50758E-02	1.50528E-02	1.50528E-02
5.0E+20	1.51035E-02	1.51035E-02	1.51033E-02	1.51030E-02	1.51024E-02	1.51005E-02	1.50975E-02	1.50917E-02	1.50757E-02	1.50528E-02
1.0E+21	1.51033E-02	1.51032E-02	1.51030E-02	1.51027E-02	1.51021E-02	1.51003E-02	1.50973E-02	1.50914E-02	1.50754E-02	1.50525E-02
2.0E+21	1.51028E-02	1.51027E-02	1.51025E-02	1.51022E-02	1.51016E-02	1.50998E-02	1.50968E-02	1.50910E-02	1.50750E-02	1.50521E-02
5.0E+21	1.51014E-02	1.51013E-02	1.51011E-02	1.51008E-02	1.51002E-02	1.50984E-02	1.50954E-02	1.50897E-02	1.50737E-02	1.50510E-02
1.0E+22	1.50991E-02	1.50991E-02	1.50989E-02	1.50986E-02	1.50980E-02	1.50962E-02	1.50932E-02	1.50917E-02	1.50491E-02	1.50491E-02
2.0E+22	1.50991E-02	1.50990E-02	1.50989E-02	1.50986E-02	1.50980E-02	1.50962E-02	1.50932E-02	1.50917E-02	1.50491E-02	1.50491E-02
5.0E+22	1.50847E-02	1.50847E-02	1.50845E-02	1.50842E-02	1.50836E-02	1.50819E-02	1.50791E-02	1.50737E-02	1.50587E-02	1.50373E-02
1.0E+23	1.50712E-02	1.50711E-02	1.50709E-02	1.50707E-02	1.50701E-02	1.50685E-02	1.50659E-02	1.50607E-02	1.50465E-02	1.50263E-02
2.0E+23	1.50543E-02	1.50543E-02	1.50541E-02	1.50538E-02	1.50533E-02	1.50519E-02	1.50516E-02	1.50447E-02	1.50316E-02	1.50131E-02
5.0E+23	1.50364E-02	1.50364E-02	1.50362E-02	1.50360E-02	1.50342E-02	1.50321E-02	1.50321E-02	1.50279E-02	1.50164E-02	1.49999E-02
1.0E+24	1.50309E-02	1.50309E-02	1.50308E-02	1.50306E-02	1.50301E-02	1.50289E-02	1.50289E-02	1.50229E-02	1.50120E-02	1.49964E-02
2.0E+24	1.50294E-02	1.50294E-02	1.50293E-02	1.50291E-02	1.50287E-02	1.50275E-02	1.50275E-02	1.50216E-02	1.50110E-02	1.49956E-02
5.0E+24	1.50289E-02	1.50289E-02	1.50288E-02	1.50286E-02	1.50282E-02	1.50270E-02	1.50270E-02	1.50212E-02	1.50106E-02	1.49954E-02
1.0E+25	1.50289E-02	1.50289E-02	1.50288E-02	1.50285E-02	1.50281E-02	1.50269E-02	1.50269E-02	1.50211E-02	1.50106E-02	1.49954E-02

Table B10. Total Photodissociation Rate Coefficients for the Process
 $\text{NO}_2 + h\nu \rightarrow \text{NO} + \text{O}^{\text{[D]}}$ for Wavelengths From 190 to 420 nm (Contd)

	03 COLUMN DENSITY				
	2.0E+17	5.0E+17	1.0E+18	2.0E+18	5.0E+18
1.0E+19	1.50170E-02	1.49479E-02	1.48792E-02	1.47941E-02	1.46491E-02
2.0E+19	1.50170E-02	1.49479E-02	1.48791E-02	1.47941E-02	1.46491E-02
5.0E+19	1.50170E-02	1.49479E-02	1.48791E-02	1.47941E-02	1.46491E-02
1.0E+20	1.49479E-02	1.48791E-02	1.48791E-02	1.47940E-02	1.46491E-02
2.0E+20	1.50169E-02	1.49478E-02	1.48791E-02	1.47940E-02	1.46491E-02
5.0E+20	1.50168E-02	1.49478E-02	1.48790E-02	1.47940E-02	1.46491E-02
1.0E+21	1.50166E-02	1.49476E-02	1.48789E-02	1.47940E-02	1.46491E-02
2.0E+21	1.50162E-02	1.49473E-02	1.48788E-02	1.47939E-02	1.46490E-02
5.0E+21	1.50153E-02	1.49467E-02	1.48784E-02	1.47937E-02	1.46490E-02
1.0E+22	1.50137E-02	1.49456E-02	1.48777E-02	1.47933E-02	1.46489E-02
2.0E+22	1.50108E-02	1.49437E-02	1.48765E-02	1.47928E-02	1.46488E-02
5.0E+22	1.50038E-02	1.49391E-02	1.48738E-02	1.47915E-02	1.46485E-02
1.0E+23	1.49947E-02	1.49333E-02	1.48704E-02	1.47899E-02	1.46482E-02
2.0E+23	1.49840E-02	1.49268E-02	1.48668E-02	1.47883E-02	1.46479E-02
5.0E+23	1.49739E-02	1.49214E-02	1.48641E-02	1.47872E-02	1.46477E-02
1.0E+24	1.49715E-02	1.49204E-02	1.48637E-02	1.47870E-02	1.46477E-02
2.0E+24	1.49711E-02	1.49204E-02	1.48637E-02	1.47870E-02	1.46476E-02
5.0E+24	1.49710E-02	1.49204E-02	1.48637E-02	1.47870E-02	1.46476E-02
1.0E+25	1.49710E-02	1.49204E-02	1.48637E-02	1.47870E-02	1.46476E-02

02 COLUMN DENSITY

Table B11. Total Photodissociation Rate Coefficients
 for the Process $\text{NO}_3 + h\nu \rightarrow \text{NO} + \text{O}_2$ for Wavelengths
 From 450 to 680 nm

1.0E+14	7.04137E-02
2.0E+14	7.04137E-02
5.0E+14	7.04136E-02
1.0E+15	7.04135E-02
2.0E+15	7.04133E-02
5.0E+15	7.04126E-02
1.0E+16	7.04115E-02
2.0E+16	7.04093E-02
5.0E+16	7.04027E-02
1.0E+17	7.03918E-02
2.0E+17	7.03699E-02
5.0E+17	7.03042E-02
1.0E+18	7.01948E-02
2.0E+18	6.99767E-02
5.0E+18	6.932271E-02
1.0E+19	6.82601E-02
2.0E+19	6.61829E-02
5.0E+19	6.03813E-02
1.0E+20	5.19884E-02
2.0E+20	3.90163E-02
5.0E+20	1.81717E-02
1.0E+21	6.74160E-03

Q3 COLUMN DENSITY

Table B12. Total Photodissociation Rate Coefficients for the Process $\text{N}_2\text{O} + h\nu \rightarrow \text{N}_2 + \text{O}^1\text{D}$ for Wavelengths From 190 to 315 nm

Q3 COLUMN DENSITY						
1.0E+11	2.0E+11	5.0E+11	1.0E+12	2.0E+12	5.0E+12	1.0E+13
2.0E+14						
1.0E+17	7.76688E-07	7.76688E-07	7.76687E-07	7.76687E-07	7.76685E-07	7.76665E-07
2.0E+17	7.76675E-07	7.76675E-07	7.76675E-07	7.76674E-07	7.76673E-07	7.76670E-07
5.0E+17	7.76637E-07	7.76637E-07	7.76636E-07	7.76636E-07	7.76634E-07	7.76632E-07
1.0E+18	7.76573E-07	7.76573E-07	7.76573E-07	7.76573E-07	7.76571E-07	7.76569E-07
2.0E+18	7.76447E-07	7.76447E-07	7.76446E-07	7.76446E-07	7.76444E-07	7.76442E-07
5.0E+18	7.76072E-07	7.76072E-07	7.76071E-07	7.76071E-07	7.76069E-07	7.76067E-07
1.0E+19	7.75460E-07	7.75460E-07	7.75460E-07	7.75460E-07	7.75458E-07	7.75456E-07
2.0E+19	7.74287E-07	7.74287E-07	7.74287E-07	7.74287E-07	7.74285E-07	7.74282E-07
5.0E+19	7.71086E-07	7.71086E-07	7.71086E-07	7.71086E-07	7.71084E-07	7.71084E-07
1.0E+20	7.66513E-07	7.66513E-07	7.66512E-07	7.66512E-07	7.66510E-07	7.66508E-07
2.0E+20	7.59021E-07	7.59021E-07	7.59021E-07	7.59020E-07	7.59019E-07	7.59017E-07
5.0E+20	7.58248E-07	7.58248E-07	7.58248E-07	7.58248E-07	7.58246E-07	7.58244E-07
1.0E+21	7.41207E-07	7.41207E-07	7.41207E-07	7.41207E-07	7.41205E-07	7.41203E-07
2.0E+21	7.20595E-07	7.20595E-07	7.20595E-07	7.20594E-07	7.20593E-07	7.20590E-07
5.0E+21	6.88458E-07	6.88458E-07	6.88458E-07	6.88458E-07	6.88456E-07	6.88454E-07
1.0E+22	6.30746E-07	6.30746E-07	6.30746E-07	6.30746E-07	6.30744E-07	6.30742E-07
2.0E+22	5.60380E-07	5.60380E-07	5.60380E-07	5.60380E-07	5.60379E-07	5.60377E-07
5.0E+22	4.41787E-07	4.41787E-07	4.41787E-07	4.41787E-07	4.41786E-07	4.41784E-07
1.0E+23	2.74667E-07	2.74667E-07	2.74667E-07	2.74667E-07	2.74665E-07	2.74663E-07
2.0E+23	1.31564E-07	1.31564E-07	1.31564E-07	1.31564E-07	1.31563E-07	1.31562E-07
5.0E+23	3.86021E-08	3.86021E-08	3.86020E-08	3.86020E-08	3.86018E-08	3.86016E-08
1.0E+24	1.89016E-08	1.89016E-08	1.89016E-08	1.89015E-08	1.89013E-08	1.89011E-08
2.0E+24	1.822305E-08	1.822305E-08	1.822305E-08	1.822305E-08	1.822301E-08	1.822296E-08
5.0E+24	1.81457E-08	1.81457E-08	1.81456E-08	1.81456E-08	1.81455E-08	1.81452E-08
1.0E+25	1.81365E-08	1.81365E-08	1.81364E-08	1.81364E-08	1.81362E-08	1.81360E-08
2.0E+25	1.81362E-08	1.81362E-08	1.81361E-08	1.81361E-08	1.81359E-08	1.81357E-08

02 COLUMN DENSITY

Table B12. Total Photodissociation Rate Coefficients for the Process $\text{N}_2\text{O} + h\nu \rightarrow \text{N}_2 + \text{O}(^1\text{D})$ for Wavelengths From 190 to 315 nm (Contd)

03 COLUMN DENSITY						
	2.0E+14	5.0E+14	1.0E+15	2.0E+15	5.0E+16	1.0E+17
1.0E+17	7.76596E-07	7.76458E-07	7.76229E-07	7.75772E-07	7.74405E-07	7.72142E-07
2.0E+17	7.76583E-07	7.76446E-07	7.76217E-07	7.75759E-07	7.74392E-07	7.72129E-07
3.0E+17	7.76545E-07	7.76407E-07	7.76178E-07	7.75721E-07	7.74354E-07	7.72091E-07
4.0E+17	7.76511E-07	7.76344E-07	7.76115E-07	7.75658E-07	7.74291E-07	7.72028E-07
5.0E+18	7.76481E-07	7.76344E-07	7.76115E-07	7.75989E-07	7.74164E-07	7.71902E-07
6.0E+18	7.76355E-07	7.76218E-07	7.75531E-07	7.75843E-07	7.75157E-07	7.71529E-07
7.0E+18	7.75980E-07	7.75843E-07	7.75614E-07	7.75157E-07	7.73790E-07	7.71529E-07
8.0E+18	7.75369E-07	7.75231E-07	7.75003E-07	7.74546E-07	7.73180E-07	7.70920E-07
9.0E+19	7.75369E-07	7.75231E-07	7.74546E-07	7.73830E-07	7.73374E-07	7.72010E-07
1.0E+19	7.74995E-07	7.74053E-07	7.73374E-07	7.73017E-07	7.69752E-07	7.65295E-07
2.0E+19	7.70995E-07	7.70858E-07	7.70630E-07	7.70175E-07	7.68816E-07	7.66566E-07
3.0E+19	7.66442E-07	7.66286E-07	7.66059E-07	7.65606E-07	7.64253E-07	7.62013E-07
4.0E+19	7.58931E-07	7.58796E-07	7.58571E-07	7.58121E-07	7.56778E-07	7.54555E-07
5.0E+19	7.58158E-07	7.58023E-07	7.57799E-07	7.57350E-07	7.56008E-07	7.53798E-07
6.0E+19	7.41119E-07	7.40986E-07	7.40765E-07	7.40324E-07	7.39005E-07	7.36824E-07
7.0E+19	7.20508E-07	7.20378E-07	7.20162E-07	7.19731E-07	7.18440E-07	7.16303E-07
8.0E+19	6.88375E-07	6.88249E-07	6.88041E-07	6.87625E-07	6.86381E-07	6.84323E-07
9.0E+19	6.30668E-07	6.30551E-07	6.30356E-07	6.29967E-07	6.28803E-07	6.26879E-07
1.0E+20	5.60309E-07	5.60202E-07	5.60024E-07	5.59669E-07	5.58608E-07	5.56855E-07
2.0E+20	4.41729E-07	4.41641E-07	4.41494E-07	4.41202E-07	4.40329E-07	4.38888E-07
3.0E+20	2.74625E-07	2.74563E-07	2.74459E-07	2.74251E-07	2.73632E-07	2.72611E-07
4.0E+20	1.31538E-07	1.31500E-07	1.31436E-07	1.31308E-07	1.30929E-07	1.30307E-07
5.0E+20	3.85889E-08	3.85692E-08	3.85364E-08	3.84712E-08	3.82778E-08	3.79634E-08
6.0E+20	1.88914E-08	1.88762E-08	1.88508E-08	1.88004E-08	1.86514E-08	1.84099E-08
7.0E+20	1.82208E-08	1.82063E-08	1.81822E-08	1.81341E-08	1.79921E-08	1.77620E-08
8.0E+20	1.81361E-08	1.81218E-08	1.80980E-08	1.80507E-08	1.79108E-08	1.76841E-08
9.0E+20	1.81269E-08	1.81127E-08	1.80889E-08	1.80417E-08	1.79020E-08	1.76756E-08
1.0E+21	1.81266E-08	1.81123E-08	1.80886E-08	1.80413E-08	1.79017E-08	1.76753E-08
2.0E+21	1.81266E-08	1.81123E-08	1.80886E-08	1.80413E-08	1.79017E-08	1.76753E-08
02 COLUMN DENSITY						
	2.0E+14	5.0E+14	1.0E+15	2.0E+15	5.0E+16	1.0E+17

Table B12. Total Photodissociation Rate Coefficients for the Process $\text{N}_2\text{O} + \text{h}\nu \rightarrow \text{N}_2 + \text{O}(1\text{D})$ for Wavelengths From 190 to 315 nm (Contd)

	02 COLUMN DENSITY					03 COLUMN DENSITY				
	5.0E+17	1.0E+18	2.0E+18	5.0E+18	1.0E+19	2.0E+19	5.0E+19	1.0E+19	2.0E+19	5.0E+19
1.0E+17	6.03763E-07	4.83551E-07	3.18465E-07	9.86375E-08	1.63090E-08	1.07089E-09	8.59453E-11			
2.0E+17	6.03753E-07	4.83543E-07	3.18460E-07	9.86362E-08	1.63089E-08	1.07089E-09	8.59453E-11			
5.0E+17	6.03723E-07	4.83519E-07	3.18444E-07	9.86322E-08	1.63084E-08	1.07088E-09	8.59453E-11			
1.0E+18	6.03672E-07	4.83479E-07	3.18419E-07	9.86255E-08	1.63077E-08	1.07087E-09	8.59453E-11			
2.0E+18	6.03572E-07	4.83399E-07	3.18368E-07	9.86123E-08	1.63062E-08	1.07085E-09	8.59453E-11			
5.0E+18	6.03273E-07	4.83161E-07	3.18217E-07	9.85730E-08	1.63018E-08	1.07078E-09	8.59453E-11			
1.0E+19	6.02787E-07	4.82774E-07	3.17971E-07	9.85087E-08	1.62946E-08	1.07066E-09	8.59453E-11			
2.0E+19	6.01853E-07	4.82030E-07	3.17497E-07	9.83847E-08	1.62806E-08	1.07044E-09	8.59453E-11			
5.0E+19	5.99303E-07	4.79995E-07	3.16199E-07	9.80421E-08	1.62413E-08	1.06979E-09	8.59453E-11			
1.0E+20	5.95653E-07	4.77078E-07	3.14330E-07	9.75427E-08	1.61828E-08	1.06877E-09	8.59453E-11			
2.0E+20	5.89662E-07	4.72279E-07	3.11242E-07	9.67042E-08	1.60817E-08	1.06692E-09	8.59452E-11			
5.0E+20	5.89092E-07	4.71844E-07	3.10977E-07	9.66372E-08	1.60738E-08	1.06678E-09	8.59452E-11			
1.0E+21	5.75467E-07	4.60915E-07	3.03913E-07	9.46837E-08	1.58296E-08	1.06198E-09	8.59451E-11			
2.0E+21	5.59003E-07	4.47700E-07	2.95339E-07	9.22722E-08	1.55172E-08	1.05539E-09	8.59449E-11			
5.0E+21	5.33511E-07	4.27298E-07	2.82119E-07	8.85198E-08	1.50175E-08	1.04420E-09	8.59446E-11			
1.0E+22	4.87436E-07	3.90206E-07	2.57815E-07	8.13909E-08	1.40144E-08	1.01943E-09	8.59437E-11			
2.0E+22	4.31625E-07	3.45375E-07	2.28419E-07	7.26317E-08	1.27358E-08	9.85761E-10	8.59425E-11			
5.0E+22	3.38947E-07	2.71525E-07	1.80395E-07	5.83839E-08	1.06352E-08	9.29097E-10	8.59403E-11			
1.0E+23	2.06846E-07	1.65135E-07	1.09845E-07	3.63482E-08	7.14610E-09	8.25112E-10	8.59359E-11			
2.0E+23	9.55816E-08	7.61533E-08	5.11481E-08	1.79497E-08	4.17913E-09	7.33719E-10	8.59318E-11			
5.0E+23	2.52507E-08	2.03050E-08	1.43418E-08	6.21662E-09	2.22475E-09	6.70637E-10	8.59289E-11			
1.0E+24	1.02400E-08	8.04676E-09	5.85509E-09	3.22418E-09	1.67355E-09	6.50942E-10	8.59279E-11			
2.0E+24	9.95878E-09	7.83253E-09	5.70855E-09	3.17142E-09	1.66353E-09	6.50572E-10	8.59278E-11			
5.0E+24	9.95734E-09	7.83247E-09	5.70855E-09	3.17142E-09	1.66353E-09	6.50572E-10	8.59278E-11			
1.0E+25	9.95724E-09	7.83247E-09	5.70855E-09	3.17142E-09	1.66353E-09	6.50572E-10	8.59278E-11			
2.0E+25	9.95723E-09	7.83247E-09	5.70855E-09	3.17142E-09	1.66353E-09	6.50572E-10	8.59278E-11			

Table B13. Total Photodissociation Rate Coefficients for the Process $\text{N}_2\text{O}_5 + h\nu \rightarrow \text{NO}_2 + \text{NO}_2 + \text{O}$ for Wavelengths From 210 to 380 nm

	03 COLUMN DENSITY									
	1.0E+11	2.0E+11	5.0E+11	1.0E+12	2.0E+12	5.0E+12	1.0E+13	2.0E+13	5.0E+13	1.0E+14
1.0E+17	6.92337E-04	6.92337E-04	6.92336E-04	6.92334E-04	6.92331E-04	6.92323E-04	6.92309E-04	6.92281E-04	6.92198E-04	6.92059E-04
2.0E+17	6.92337E-04	6.92336E-04	6.92335E-04	6.92334E-04	6.92331E-04	6.92323E-04	6.92309E-04	6.92281E-04	6.92198E-04	6.92059E-04
5.0E+17	6.92336E-04	6.92336E-04	6.92335E-04	6.92334E-04	6.92331E-04	6.92322E-04	6.92308E-04	6.92280E-04	6.92197E-04	6.92058E-04
1.0E+18	6.92335E-04	6.92334E-04	6.92334E-04	6.92332E-04	6.92329E-04	6.92321E-04	6.92307E-04	6.92279E-04	6.92196E-04	6.92056E-04
2.0E+18	6.92332E-04	6.92332E-04	6.92331E-04	6.92330E-04	6.92327E-04	6.92321E-04	6.92277E-04	6.92277E-04	6.92193E-04	6.92054E-04
5.0E+18	6.92326E-04	6.92325E-04	6.92324E-04	6.92323E-04	6.92320E-04	6.92312E-04	6.92298E-04	6.92270E-04	6.92187E-04	6.92047E-04
1.0E+19	6.92314E-04	6.92314E-04	6.92313E-04	6.92312E-04	6.92310E-04	6.92309E-04	6.92287E-04	6.92259E-04	6.92175E-04	6.92036E-04
2.0E+19	6.92291E-04	6.92290E-04	6.92289E-04	6.92286E-04	6.92278E-04	6.92278E-04	6.92264E-04	6.92236E-04	6.92152E-04	6.92013E-04
5.0E+19	6.92223E-04	6.92222E-04	6.92222E-04	6.92220E-04	6.92217E-04	6.92209E-04	6.92195E-04	6.92167E-04	6.92084E-04	6.91944E-04
1.0E+20	6.92108E-04	6.92108E-04	6.92107E-04	6.92106E-04	6.92103E-04	6.92095E-04	6.92081E-04	6.92053E-04	6.91969E-04	6.91830E-04
2.0E+20	6.91880E-04	6.91880E-04	6.91879E-04	6.91879E-04	6.91875E-04	6.91866E-04	6.91852E-04	6.91824E-04	6.91741E-04	6.91602E-04
5.0E+20	6.91195E-04	6.91195E-04	6.91194E-04	6.91193E-04	6.91190E-04	6.91182E-04	6.91168E-04	6.91140E-04	6.91057E-04	6.90917E-04
1.0E+21	6.90058E-04	6.90057E-04	6.90057E-04	6.90056E-04	6.90052E-04	6.90044E-04	6.90030E-04	6.9002E-04	6.89912E-04	6.89780E-04
2.0E+21	6.87794E-04	6.87793E-04	6.87793E-04	6.87791E-04	6.87788E-04	6.87780E-04	6.87766E-04	6.87738E-04	6.87716E-04	6.87755E-04
5.0E+21	6.81093E-04	6.81093E-04	6.81092E-04	6.81090E-04	6.81088E-04	6.81079E-04	6.81066E-04	6.81038E-04	6.80955E-04	6.80817E-04
1.0E+22	6.70223E-04	6.70222E-04	6.70220E-04	6.70217E-04	6.70217E-04	6.70209E-04	6.70195E-04	6.70168E-04	6.70086E-04	6.69949E-04
2.0E+22	6.49550E-04	6.49550E-04	6.49549E-04	6.49548E-04	6.49545E-04	6.49537E-04	6.49524E-04	6.49497E-04	6.49461E-04	6.49281E-04
5.0E+22	5.95214E-04	5.95214E-04	5.95213E-04	5.95212E-04	5.95210E-04	5.95202E-04	5.95189E-04	5.95163E-04	5.95086E-04	5.94957E-04
1.0E+23	5.25402E-04	5.25402E-04	5.25401E-04	5.25400E-04	5.25398E-04	5.25390E-04	5.25354E-04	5.25281E-04	5.25160E-04	5.25100E-04
2.0E+23	4.37587E-04	4.37587E-04	4.37586E-04	4.37585E-04	4.37583E-04	4.37576E-04	4.37565E-04	4.37533E-04	4.37477E-04	4.37367E-04
5.0E+23	3.45059E-04	3.45059E-04	3.45058E-04	3.45058E-04	3.45056E-04	3.45050E-04	3.45041E-04	3.45022E-04	3.44965E-04	3.44870E-04
1.0E+24	3.16513E-04	3.16512E-04	3.16511E-04	3.16511E-04	3.16509E-04	3.16504E-04	3.16478E-04	3.16425E-04	3.16338E-04	3.16338E-04
2.0E+24	3.06849E-04	3.06849E-04	3.06848E-04	3.06847E-04	3.06846E-04	3.06840E-04	3.06832E-04	3.06815E-04	3.06681E-04	3.06656E-04
5.0E+24	3.02786E-04	3.02785E-04	3.02784E-04	3.02784E-04	3.02783E-04	3.02778E-04	3.02769E-04	3.02753E-04	3.02704E-04	3.02622E-04
1.0E+25	3.02265E-04	3.02285E-04	3.02284E-04	3.02283E-04	3.02282E-04	3.02277E-04	3.02269E-04	3.02252E-04	3.02212E-04	3.02104E-04
2.0E+25	3.02267E-04	3.02267E-04	3.02266E-04	3.02266E-04	3.02259E-04	3.02251E-04	3.02235E-04	3.022104E-04	3.02185E-04	3.02104E-04

Table B13. Total Photodissociation Rate Coefficients for the Process $\text{N}_2\text{C}_5 + h\nu \rightarrow \text{NO}_2 + \text{NO}_2 + \text{O}$ for Wavelengths From 210 to 380 nm (Contd)

03 COLUMN DENSITY						
2.0E+14	5.0E+14	1.0E+15	2.0E+15	5.0E+15	1.0E+16	2.0E+16
1.0E+17	6.91780E-04	6.90946E-04	6.89561E-04	6.86805E-04	6.78656E-04	7.72142E-07
2.0E+17	6.91780E-04	6.90946E-04	6.89561E-04	6.86805E-04	6.78656E-04	7.67675E-07
3.0E+17	6.91779E-04	6.90946E-04	6.89560E-04	6.86804E-04	6.78655E-04	7.72129E-07
4.0E+17	6.91778E-04	6.90944E-04	6.89559E-04	6.86803E-04	6.78654E-04	7.72091E-07
5.0E+17	6.91778E-04	6.90944E-04	6.89559E-04	6.86803E-04	6.78654E-04	7.67624E-07
1.0E+18	6.91776E-04	6.90942E-04	6.89557E-04	6.86801E-04	6.78652E-04	7.72028E-07
2.0E+18	6.91776E-04	6.90942E-04	6.89557E-04	6.86801E-04	6.78652E-04	7.67561E-07
3.0E+18	6.91776E-04	6.90942E-04	6.89557E-04	6.86801E-04	6.78652E-04	7.67436E-07
4.0E+18	6.91769E-04	6.90935E-04	6.89550E-04	6.86794E-04	6.78645E-04	7.71529E-07
5.0E+18	6.91758E-04	6.90924E-04	6.89538E-04	6.86782E-04	6.78634E-04	7.70920E-07
6.0E+19	6.91735E-04	6.90911E-04	6.89516E-04	6.86760E-04	6.78611E-04	7.69752E-07
7.0E+19	6.91666E-04	6.90833E-04	6.89447E-04	6.86691E-04	6.78543E-04	7.66566E-07
8.0E+19	6.91552E-04	6.90718E-04	6.89333E-04	6.86577E-04	6.78430E-04	7.62013E-07
9.0E+19	6.91324E-04	6.90490E-04	6.89105E-04	6.86350E-04	6.78204E-04	7.54555E-07
1.0E+20	6.90639E-04	6.89806E-04	6.88422E-04	6.85668E-04	6.77526E-04	7.53788E-07
2.0E+20	6.90639E-04	6.89806E-04	6.88422E-04	6.85668E-04	6.77526E-04	7.49405E-07
3.0E+20	6.89502E-04	6.88726E-04	6.87286E-04	6.84535E-04	6.76400E-04	7.36824E-07
4.0E+20	6.89502E-04	6.88726E-04	6.87286E-04	6.84535E-04	6.76400E-04	7.32518E-07
5.0E+20	6.87239E-04	6.86408E-04	6.85027E-04	6.82280E-04	6.74158E-04	7.16305E-07
6.0E+20	6.80541E-04	6.79714E-04	6.78340E-04	6.75607E-04	6.67525E-04	6.45925E-07
7.0E+20	6.69675E-04	6.68855E-04	6.67492E-04	6.64781E-04	6.56765E-04	6.50168E-07
8.0E+20	6.49011E-04	6.48204E-04	6.46862E-04	6.44194E-04	6.36305E-04	5.56855E-07
9.0E+20	5.94699E-04	5.93926E-04	5.92642E-04	5.90089E-04	5.82543E-04	4.38888E-07
1.0E+21	5.24918E-04	5.24193E-04	5.22987E-04	5.20591E-04	5.13514E-04	2.72611E-07
2.0E+21	4.37147E-04	4.36488E-04	4.35394E-04	4.33218E-04	4.26798E-04	1.30307E-07
3.0E+21	3.44680E-04	3.44112E-04	3.43169E-04	3.41296E-04	3.35771E-04	1.29099E-07
4.0E+21	3.16163E-04	3.15639E-04	3.14769E-04	3.13042E-04	3.07949E-04	1.84099E-08
5.0E+21	3.06513E-04	3.06011E-04	3.05177E-04	3.03521E-04	2.988638E-04	1.77620E-08
6.0E+21	3.02457E-04	3.01966E-04	3.01150E-04	2.99528E-04	2.94749E-04	1.76841E-08
7.0E+21	3.01958E-04	3.01467E-04	3.00653E-04	2.99036E-04	2.94271E-04	1.76756E-08
8.0E+21	3.01940E-04	3.01450E-04	3.00636E-04	2.99019E-04	2.94254E-04	1.76753E-08
02 COLUMN DENSITY						
2.0E+14	5.0E+14	1.0E+15	2.0E+15	5.0E+15	1.0E+16	2.0E+16

Table B13. Total Photodissociation Rate Coefficients for the Process
 $\text{N}_2\text{O}_5 + h\nu \rightarrow \text{NO}_2 + \text{NO}_2 + \text{O}$ for Wavelengths From 210 to 380 nm (Contd)

	O3 COLUMN DENSITY				
	5.0E+17	1.0E+18	2.0E+18	5.0E+18	1.0E+19
1.0E+17	6.03763E-07	4.83551E-07	3.18465E-07	9.86375E-08	1.63090E-08
2.0E+17	6.03753E-07	4.83543E-07	3.18460E-07	9.86362E-08	1.63089E-08
5.0E+17	6.03723E-07	4.83519E-07	3.18444E-07	9.86322E-08	1.63084E-08
1.0E+18	6.03672E-07	4.83479E-07	3.18419E-07	9.86255E-08	1.63077E-08
2.0E+18	6.03572E-07	4.83399E-07	3.18368E-07	9.86123E-08	1.63062E-08
5.0E+18	6.03273E-07	4.83161E-07	3.18217E-07	9.85730E-08	1.63018E-08
1.0E+19	6.02787E-07	4.82774E-07	3.17971E-07	9.85087E-08	1.62946E-08
2.0E+19	6.01853E-07	4.82030E-07	3.17497E-07	9.83847E-08	1.62806E-08
5.0E+19	5.99303E-07	4.79995E-07	3.16199E-07	9.80421E-08	1.62413E-08
1.0E+20	5.95633E-07	4.77078E-07	3.14330E-07	9.75427E-08	1.61828E-08
2.0E+20	5.89966E-07	4.72279E-07	3.11242E-07	9.67042E-08	1.60817E-08
5.0E+20	5.80922E-07	4.71644E-07	3.10977E-07	9.66372E-08	1.60738E-08
1.0E+21	5.75467E-07	4.60915E-07	3.10913E-07	9.46837E-08	1.58296E-08
2.0E+21	5.59003E-07	4.47700E-07	2.95339E-07	9.22722E-08	1.55172E-08
5.0E+21	5.33511E-07	4.27298E-07	2.82119E-07	8.85198E-08	1.50175E-08
1.0E+22	4.87436E-07	3.90206E-07	2.57815E-07	8.13909E-08	1.40144E-08
2.0E+22	4.31625E-07	3.45375E-07	2.28419E-07	7.26317E-08	1.27358E-08
5.0E+22	3.3897E-07	2.71525E-07	1.80395E-07	5.83839E-08	1.06352E-08
1.0E+23	2.06846E-07	1.65135E-07	1.09845E-07	3.63482E-08	7.14610E-09
2.0E+23	9.55816E-08	7.61533E-08	5.11481E-08	1.79497E-08	4.17913E-09
5.0E+23	2.52507E-08	2.03050E-08	1.43418E-08	6.21662E-09	2.22475E-09
1.0E+24	1.02400E-08	8.04676E-09	5.85509E-09	3.22418E-09	1.67355E-09
2.0E+24	9.95879E-09	7.83253E-09	5.70855E-09	3.17142E-09	1.66353E-09
5.0E+24	9.95734E-09	7.83247E-09	5.70855E-09	3.17142E-09	1.66353E-09
1.0E+25	9.95724E-09	7.83247E-09	5.70855E-09	3.17142E-09	1.66353E-09
2.0E+25	9.95723E-09	7.83247E-09	5.70855E-09	3.17142E-09	1.66353E-09

Table B14. Total Photodissociation Rate Coefficients for the Process $\text{HNO}_2 + h\nu \rightarrow \text{OH} + \text{NO}$ for Wavelengths From 300 to 390 nm

	1.0E+14	5.84756E-04	Q3 COLUMN DENSITY
2.0E+14	5.84755E-04		
5.0E+14	5.84754E-04		
1.0E+15	5.84753E-04		
2.0E+15	5.84749E-04		
5.0E+15	5.84739E-04		
1.0E+16	5.84722E-04		
2.0E+16	5.84688E-04		
5.0E+16	5.84585E-04		
1.0E+17	5.84416E-04		
2.0E+17	5.84083E-04		
5.0E+17	5.83115E-04		
1.0E+18	5.81605E-04		
2.0E+18	5.78914E-04		
5.0E+18	5.72656E-04		
1.0E+19	5.65442E-04		
2.0E+19	5.55729E-04		
5.0E+19	5.37947E-04		
1.0E+20	5.20145E-04		
2.0E+20	4.99103E-04		
5.0E+20	4.68331E-04		
1.0E+21	4.43897E-04		

Table B15. Total Photodissociation Rate Coefficients for the Process $\text{HNO}_3 + h\nu \rightarrow \text{OH} + \text{NO}_2$ for Wavelengths From 192 to 325 nm

	O3 COLUMN DENSITY									
	1.0E+12	2.0E+12	5.0E+12	1.0E+13	2.0E+13	5.0E+13	1.0E+14	2.0E+14	5.0E+14	1.0E+15
1.0E+17	1.16539E-04	1.16539E-04	1.16539E-04	1.16539E-04	1.16537E-04	1.16535E-04	1.16528E-04	1.16493E-04	1.16493E-04	1.16424E-04
2.0E+17	1.16539E-04	1.16539E-04	1.16538E-04	1.16537E-04	1.16535E-04	1.16528E-04	1.16516E-04	1.16493E-04	1.16493E-04	1.16424E-04
5.0E+17	1.16539E-04	1.16539E-04	1.16538E-04	1.16537E-04	1.16534E-04	1.16528E-04	1.16516E-04	1.16493E-04	1.16493E-04	1.16424E-04
1.0E+18	1.16538E-04	1.16538E-04	1.16537E-04	1.16536E-04	1.16534E-04	1.16527E-04	1.16515E-04	1.16492E-04	1.16492E-04	1.16423E-04
2.0E+18	1.16537E-04	1.16537E-04	1.16536E-04	1.16535E-04	1.16532E-04	1.16525E-04	1.16514E-04	1.16491E-04	1.16491E-04	1.16421E-04
5.0E+18	1.16532E-04	1.16532E-04	1.16530E-04	1.16530E-04	1.16528E-04	1.16521E-04	1.16510E-04	1.16479E-04	1.16479E-04	1.16417E-04
1.0E+19	1.16525E-04	1.16525E-04	1.16524E-04	1.16523E-04	1.16521E-04	1.16514E-04	1.16502E-04	1.16479E-04	1.16479E-04	1.16410E-04
2.0E+19	1.16511E-04	1.16511E-04	1.16509E-04	1.16509E-04	1.16507E-04	1.16500E-04	1.16488E-04	1.16465E-04	1.16465E-04	1.16280E-04
5.0E+19	1.16469E-04	1.16469E-04	1.16468E-04	1.16467E-04	1.16465E-04	1.16458E-04	1.16446E-04	1.16434E-04	1.16434E-04	1.16238E-04
1.0E+20	1.16400E-04	1.16400E-04	1.16399E-04	1.16398E-04	1.16396E-04	1.16389E-04	1.16377E-04	1.16354E-04	1.16354E-04	1.16169E-04
2.0E+20	1.16265E-04	1.16265E-04	1.16264E-04	1.16263E-04	1.16261E-04	1.16254E-04	1.16242E-04	1.16219E-04	1.16219E-04	1.16035E-04
5.0E+20	1.16127E-04	1.16127E-04	1.16126E-04	1.16125E-04	1.16122E-04	1.16116E-04	1.16104E-04	1.16081E-04	1.16081E-04	1.16012E-04
1.0E+21	1.15652E-04	1.15652E-04	1.15652E-04	1.15650E-04	1.15648E-04	1.15641E-04	1.15630E-04	1.15606E-04	1.15537E-04	1.15422E-04
2.0E+21	1.14835E-04	1.14835E-04	1.14834E-04	1.14833E-04	1.14831E-04	1.14824E-04	1.14812E-04	1.14789E-04	1.14789E-04	1.14605E-04
5.0E+21	1.12858E-04	1.12857E-04	1.12857E-04	1.12856E-04	1.12853E-04	1.12846E-04	1.12835E-04	1.12812E-04	1.12743E-04	1.12629E-04
1.0E+22	1.09086E-04	1.09085E-04	1.09085E-04	1.09084E-04	1.09081E-04	1.09075E-04	1.09063E-04	1.09041E-04	1.08860E-04	1.08860E-04

Table B15. Total Photodissociation Rate Coefficients for the Process $\text{HNO}_3 + h\nu \rightarrow \text{OH} + \text{NO}_2$ for Wavelengths From 192 to 325 nm (Contd)

		O3 COLUMN DENSITY									
		2.0E+15	5.0E+15	1.0E+16	2.0E+16	5.0E+16	1.0E+17	2.0E+17	5.0E+17	1.0E+18	2.0E+18
1.0E+17	1.16079E-04	1.15398E-04	1.14292E-04	1.12176E-04	1.06523E-04	9.89202E-05	8.80795E-05	6.98619E-05	5.31722E-05	3.41525E-05	
2.0E+17	1.16079E-04	1.15398E-04	1.14292E-04	1.12176E-04	1.06523E-04	9.89201E-05	8.80794E-05	6.98618E-05	5.31721E-05	3.41525E-05	
5.0E+17	1.16078E-04	1.15398E-04	1.14291E-04	1.12176E-04	1.06522E-04	9.89197E-05	8.80790E-05	6.98614E-05	5.31718E-05	3.41523E-05	
1.0E+18	1.16078E-04	1.15397E-04	1.14290E-04	1.12175E-04	1.06522E-04	9.89190E-05	8.80784E-05	6.98609E-05	5.31713E-05	3.41520E-05	
2.0E+18	1.16076E-04	1.15396E-04	1.14289E-04	1.12174E-04	1.06520E-04	9.89177E-05	8.80771E-05	6.98597E-05	5.31704E-05	3.41514E-05	
5.0E+18	1.16072E-04	1.15391E-04	1.14285E-04	1.12169E-04	1.06516E-04	9.89136E-05	8.80733E-05	6.98564E-05	5.31677E-05	3.41495E-05	
1.0E+19	1.16065E-04	1.15384E-04	1.14278E-04	1.12162E-04	1.06509E-04	9.89069E-05	8.80669E-05	6.98508E-05	5.31631E-05	3.41464E-05	
2.0E+19	1.16051E-04	1.15370E-04	1.14264E-04	1.12148E-04	1.06495E-04	9.88935E-05	8.80541E-05	6.98396E-05	5.31540E-05	3.41402E-05	
5.0E+19	1.16009E-04	1.15328E-04	1.14222E-04	1.12107E-04	1.06454E-04	9.88536E-05	8.80161E-05	6.98064E-05	5.31269E-05	3.41217E-05	
1.0E+20	1.15940E-04	1.15260E-04	1.14153E-04	1.12038E-04	1.06387E-04	9.87880E-05	8.79536E-05	6.97518E-05	5.30825E-05	3.40914E-05	
2.0E+20	1.15805E-04	1.15125E-04	1.14019E-04	1.11905E-04	1.06256E-04	9.86602E-05	8.78319E-05	6.96455E-05	5.29960E-05	3.40325E-05	
5.0E+20	1.15667E-04	1.14987E-04	1.13882E-04	1.11795E-04	1.06124E-04	9.85336E-05	8.77152E-05	6.95513E-05	5.29266E-05	3.39918E-05	
1.0E+21	1.15193E-04	1.14514E-04	1.13411E-04	1.11301E-04	1.05664E-04	9.80871E-05	8.72930E-05	6.91883E-05	5.26366E-05	3.37992E-05	
2.0E+21	1.14377E-04	1.13700E-04	1.12599E-04	1.10494E-04	1.04873E-04	9.73209E-05	8.65711E-05	6.85728E-05	5.21498E-05	3.34803E-05	
5.0E+21	1.12402E-04	1.11730E-04	1.06366E-04	1.08546E-04	1.02968E-04	9.54811E-05	8.48483E-05	6.71253E-05	5.10251E-05	3.27621E-05	
1.0E+22	1.08636E-04	1.07971E-04	1.06890E-04	1.04826E-04	9.93215E-05	9.19509E-05	8.15257E-05	6.42989E-05	4.87946E-05	3.13045E-05	

Table B20. Total Photodissociation Rate Coefficients for the Process $\text{HO}_2 + \text{h}\nu \rightarrow \text{O}_2 + \text{H}$ for Wavelengths From 180 to 274 nm

03 COLUMN DENSITY						
	2.0E+15	5.0E+15	1.0E+16	2.0E+16	5.0E+16	1.0E+17
1.0E+17	1.12764E-03	1.11447E-03	1.09309E-03	1.05241E-03	9.45069E-04	8.04875E-04
2.0E+17	1.12762E-03	1.11445E-03	1.09307E-03	1.05239E-03	9.45050E-04	8.04857E-04
5.0E+17	1.12757E-03	1.11439E-03	1.05302E-03	1.05234E-03	9.44997E-04	8.04805E-04
1.0E+18	1.12748E-03	1.11431E-03	1.09253E-03	1.05226E-03	9.44916E-04	8.04728E-04
2.0E+18	1.12734E-03	1.11417E-03	1.49279E-03	1.05212E-03	9.44773E-04	8.04595E-04
5.0E+18	1.12730E-03	1.11384E-03	1.05179E-03	1.05179E-03	9.44463E-04	8.04292E-04
1.0E+19	1.12662E-03	1.11344E-03	1.09207E-03	1.05140E-03	9.44081E-04	8.03924E-04
2.0E+19	1.12664E-03	1.11387E-03	1.09150E-03	1.05083E-03	9.43523E-04	8.03387E-04
5.0E+19	1.12487E-03	1.11170E-03	1.09034E-03	1.04968E-03	9.42404E-04	8.02312E-04
1.0E+20	1.12355E-03	1.11039E-03	1.68903E-03	1.04838E-03	9.41138E-04	8.01099E-04
2.0E+20	1.12165E-03	1.10843E-03	1.08714E-03	1.04652E-03	9.39327E-04	7.99371E-04
5.0E+20	1.12007E-03	1.10692E-03	1.06559E-03	1.04500E-03	9.37898E-04	7.98076E-04
1.0E+21	1.11508E-03	1.10195E-03	1.08066E-03	1.04013E-03	9.33225E-04	7.93687E-04
2.0E+21	1.10762E-03	1.09454E-03	1.07330E-03	1.03290E-03	9.26328E-04	7.87288E-04
5.0E+21	1.08948E-03	1.07650E-03	1.05544E-03	1.01537E-03	9.09742E-04	7.72075E-04
1.0E+22	1.05983E-03	1.04702E-03	1.02624E-03	9.86722E-04	8.82631E-04	7.47200E-04
2.0E+22	1.007345E-03	9.94850E-04	9.7446C5E-04	9.36127E-04	8.34937E-04	7.03715E-04
5.0E+22	8.75153E-04	6.63526E-04	8.44722E-04	8.09046E-04	7.15656E-04	5.95698E-04
1.0E+23	7.01310E-04	6.50916E-04	6.74107E-04	6.42298E-04	5.59643E-04	4.55085E-04
2.0E+23	4.85107E-04	4.76452E-04	4.62485E-04	4.36185E-04	3.68712E-04	2.85655E-04
5.0E+23	2.50555E-04	2.44359E-04	2.34434E-04	2.15892E-04	1.69502E-04	1.15444E-04
1.0E+24	1.70615E-04	1.65677E-04	1.57771E-04	1.43098E-04	1.66935E-04	1.43098E-04
2.0E+24	1.42977E-04	1.38649E-04	1.31728E-04	1.18914E-04	8.75310E-05	5.26373E-05
5.0E+24	1.31568E-04	1.27536E-04	1.21090E-04	1.09166E-04	8.00283E-05	4.77828E-05
1.0E+25	1.30174E-04	1.26179E-04	1.19793E-04	1.07980E-04	7.91231E-05	4.72054E-05
2.0E+25	1.30125E-04	1.26131E-04	1.19747E-04	1.07939E-04	7.90911E-05	4.71850E-05
5.0E+25						

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Table B20. Total Photodissociation Rate Coefficients for the I₁ process HO₂ + hν → O₂ + H for Wavelengths From 180 to 274 nm

	03 COLUMN DENSITY									
	1.0E+12	2.0E+12	5.0E+12	1.0E+13	2.0E+13	5.0E+13	1.0E+14	2.0E+14	5.0E+14	1.0E+15
1.0E+17	1.13657E-03	1.13656E-03	1.13655E-03	1.13653E-03	1.13648E-03	1.13635E-03	1.13612E-03	1.13567E-03	1.13433E-03	1.13209E-03
2.0E+17	1.13655E-03	1.13654E-03	1.13651E-03	1.13651E-03	1.13646E-03	1.13633E-03	1.13610E-03	1.13565E-03	1.13431E-03	1.13207E-03
5.0E+17	1.13649E-03	1.13649E-03	1.13645E-03	1.13645E-03	1.13641E-03	1.13627E-03	1.13605E-03	1.13560E-03	1.13425E-03	1.13202E-03
1.0E+18	1.13641E-03	1.13640E-03	1.13639E-03	1.13639E-03	1.13637E-03	1.13632E-03	1.13619E-03	1.13596E-03	1.13417E-03	1.13193E-03
2.0E+18	1.13627E-03	1.13626E-03	1.13625E-03	1.13625E-03	1.13623E-03	1.13618E-03	1.13605E-03	1.13537E-03	1.13403E-03	1.13179E-03
5.0E+18	1.13594E-03	1.13593E-03	1.13592E-03	1.13590E-03	1.13585E-03	1.13572E-03	1.13549E-03	1.13504E-03	1.13370E-03	1.13146E-03
1.0E+19	1.13554E-03	1.13554E-03	1.13552E-03	1.13550E-03	1.13546E-03	1.13532E-03	1.13510E-03	1.13465E-03	1.13330E-03	1.13107E-03
2.0E+19	1.13499E-03	1.13496E-03	1.13494E-03	1.13492E-03	1.13488E-03	1.13474E-03	1.13407E-03	1.13452E-03	1.13272E-03	1.13049E-03
5.0E+19	1.13379E-03	1.13378E-03	1.13379E-03	1.13378E-03	1.13377E-03	1.13371E-03	1.13355E-03	1.13290E-03	1.13156E-03	1.12932E-03
1.0E+20	1.13247E-03	1.13246E-03	1.13245E-03	1.13243E-03	1.13242E-03	1.13225E-03	1.13202E-03	1.13158E-03	1.13023E-03	1.12800E-03
2.0E+20	1.13056E-03	1.13056E-03	1.13055E-03	1.13052E-03	1.13048E-03	1.13034E-03	1.13012E-03	1.12967E-03	1.12833E-03	1.12609E-03
5.0E+20	1.12893E-03	1.12897E-03	1.12896E-03	1.12894E-03	1.12889E-03	1.12876E-03	1.12853E-03	1.12808E-03	1.12674E-03	1.12451E-03
1.0E+21	1.12397E-03	1.12396E-03	1.12396E-03	1.12395E-03	1.12394E-03	1.12376E-03	1.12353E-03	1.12308E-03	1.12174E-03	1.11951E-03
2.0E+21	1.11649E-03	1.11649E-03	1.11648E-03	1.11648E-03	1.11641E-03	1.11628E-03	1.11605E-03	1.11561E-03	1.11427E-03	1.11205E-03
5.0E+21	1.09828E-03	1.09827E-03	1.09826E-03	1.09824E-03	1.09824E-03	1.09819E-03	1.09816E-03	1.09784E-03	1.09740E-03	1.09587E-03
1.0E+22	1.06851E-03	1.06851E-03	1.06850E-03	1.06847E-03	1.06847E-03	1.06843E-03	1.06830E-03	1.06764E-03	1.06633E-03	1.06416E-03
2.0E+22	1.01580E-03	1.01580E-03	1.01578E-03	1.01576E-03	1.01572E-03	1.01559E-03	1.01538E-03	1.01495E-03	1.01368E-03	1.01156E-03
5.0E+22	8.83007E-04	8.83003E-04	8.82991E-04	8.82971E-04	8.82932E-04	8.827613E-04	8.82218E-04	8.81031E-04	8.79058E-04	8.77044E-04
1.0E+23	7.08365E-04	7.08363E-04	7.08352E-04	7.08334E-04	7.08299E-04	7.08192E-04	7.08014E-04	7.07659E-04	7.04827E-04	7.04827E-04
2.0E+23	4.90999E-04	4.90990E-04	4.90981E-04	4.90967E-04	4.90954E-04	4.90937E-04	4.90914E-04	4.90699E-04	4.90403E-04	4.88040E-04
5.0E+23	2.54775E-04	2.54772E-04	2.54766E-04	2.54755E-04	2.54734E-04	2.54670E-04	2.54351E-04	2.53714E-04	2.52656E-04	2.52294E-04
1.0E+24	1.73989E-04	1.73987E-04	1.73982E-04	1.73974E-04	1.73957E-04	1.73905E-04	1.73650E-04	1.73140E-04	1.72294E-04	1.72294E-04
2.0E+24	1.45934E-04	1.45934E-04	1.45930E-04	1.45922E-04	1.45907E-04	1.45862E-04	1.45788E-04	1.45638E-04	1.45191E-04	1.44449E-04
5.0E+24	1.34326E-04	1.34325E-04	1.34321E-04	1.34314E-04	1.34300E-04	1.34258E-04	1.34188E-04	1.34049E-04	1.33632E-04	1.32941E-04
1.0E+25	1.32907E-04	1.32906E-04	1.32901E-04	1.32894E-04	1.32881E-04	1.32839E-04	1.32770E-04	1.32632E-04	1.32219E-04	1.31534E-04
2.0E+25	1.32857E-04	1.32855E-04	1.32851E-04	1.32844E-04	1.32830E-04	1.32779E-04	1.32729E-04	1.32582E-04	1.32169E-04	1.31484E-04

Table B19. Total Photodissociation Rate Coefficients for the Process $\text{CH}_2\text{O} + h\nu \rightarrow \text{CHO} + \text{H}$ and $\text{CH}_2\text{O} + h\nu \rightarrow \text{CO} + \text{H}_2$ for Wavelengths From 240 to 360 nm (Contd)

03 COLUMN DENSITY						
5.0E+15	1.0E+16	2.0E+16	5.0E+16	1.0E+17	2.0E+17	5.0E+18
1.0E+13	2.58223E-04	2.57561E-04	2.56282E-04	2.52774E-04	2.47810E-04	2.40115E-04
2.0E+13	2.58223E-04	2.57561E-04	2.56282E-04	2.52774E-04	2.47810E-04	2.40115E-04
5.0E+13	2.58223E-04	2.57561E-04	2.56282E-04	2.52774E-04	2.47810E-04	2.40115E-04
1.0E+14	2.58222E-04	2.57560E-04	2.56281E-04	2.52773E-04	2.47809E-04	2.40114E-04
2.0E+14	2.58221E-04	2.57559E-04	2.56280E-04	2.52772E-04	2.47809E-04	2.40114E-04
5.0E+14	2.58219E-04	2.57556E-04	2.56278E-04	2.52770E-04	2.47806E-04	2.40112E-04
1.0E+15	2.58214E-04	2.57552E-04	2.56273E-04	2.52766E-04	2.47803E-04	2.40110E-04
2.0E+15	2.58205E-04	2.57543E-04	2.56265E-04	2.52758E-04	2.47796E-04	2.40104E-04
5.0E+15	2.58178E-04	2.57517E-04	2.56239E-04	2.52735E-04	2.47776E-04	2.40089E-04
1.0E+16	2.58135E-04	2.57474E-04	2.56198E-04	2.52697E-04	2.47743E-04	2.40064E-04
2.0E+16	2.58054E-04	2.57295E-04	2.56121E-04	2.52626E-04	2.47682E-04	2.40017E-04
5.0E+16	2.57850E-04	2.57193E-04	2.55925E-04	2.52446E-04	2.47525E-04	2.39895E-04
1.0E+17	2.57602E-04	2.56948E-04	2.55687E-04	2.52225E-04	2.47329E-04	2.39739E-04
2.0E+17	2.57296E-04	2.56645E-04	2.55390E-04	2.51946E-04	2.47076E-04	2.39529E-04
5.0E+17	2.56836E-04	2.56189E-04	2.54939E-04	2.51511E-04	2.46666E-04	2.39163E-04
1.0E+18	2.56368E-04	2.55722E-04	2.54476E-04	2.51060E-04	2.46234E-04	2.38765E-04
2.0E+18	2.55808E-04	2.55164E-04	2.53923E-04	2.50520E-04	2.45716E-04	2.38288E-04
5.0E+18	2.55285E-04	2.54644E-04	2.53407E-04	2.50443E-04	2.45233E-04	2.37844E-04
1.0E+19	2.55180E-04	2.54544E-04	2.53303E-04	2.49916E-04	2.45136E-04	2.37755E-04
2.0E+19	2.55168E-04	2.54528E-04	2.53291E-04	2.49904E-04	2.45125E-04	2.37744E-04
5.0E+19	2.55169E-04	2.54527E-04	2.53291E-04	2.49904E-04	2.45125E-04	2.37744E-04
1.0E+20	2.55163E-04	2.54527E-04	2.53291E-04	2.49904E-04	2.45125E-04	2.37744E-04
2.0E+20	2.55168E-04	2.54527E-04	2.53290E-04	2.49904E-04	2.45124E-04	2.37744E-04
5.0E+20	2.55168E-04	2.54527E-04	2.53290E-04	2.49904E-04	2.45125E-04	2.37744E-04
1.0E+21	2.55168E-04	2.54527E-04	2.53291E-04	2.49904E-04	2.45125E-04	2.37744E-04
2.0E+21	2.55168E-04	2.54527E-04	2.53291E-04	2.49904E-04	2.45125E-04	2.37744E-04
5.0E+21	2.55168E-04	2.54527E-04	2.53291E-04	2.49904E-04	2.45124E-04	2.37744E-04
1.0E+22	2.55167E-04	2.54527E-04	2.53291E-04	2.49904E-04	2.45124E-04	2.37744E-04
2.0E+22	2.55167E-04	2.54526E-04	2.53290E-04	2.49903E-04	2.45125E-04	2.37744E-04
5.0E+22	2.55167E-04	2.54526E-04	2.53290E-04	2.49903E-04	2.45125E-04	2.37744E-04
1.0E+23	2.55167E-04	2.54526E-04	2.53288E-04	2.49902E-04	2.45123E-04	2.37744E-04
2.0E+23	2.55167E-04	2.54526E-04	2.53288E-04	2.49902E-04	2.45122E-04	2.37743E-04
5.0E+23	2.55157E-04	2.54527E-04	2.53281E-04	2.49904E-04	2.45124E-04	2.37743E-04
1.0E+24	2.55127E-04	2.54447E-04	2.53218E-04	2.49848E-04	2.45089E-04	2.37730E-04
2.0E+24	2.55087E-04	2.54450E-04	2.53220E-04	2.49850E-04	2.45090E-04	2.37730E-04
5.0E+24	2.55084E-04	2.54447E-04	2.53218E-04	2.49848E-04	2.45089E-04	2.37730E-04
1.0E+25	2.55084E-04	2.54447E-04	2.53218E-04	2.49848E-04	2.45089E-04	2.37730E-04

02 COLUMN DENSITY

Table B11. Total Photodissociation Rate Coefficients for the Process $\text{CH}_2\text{O} + \text{h}\nu \rightarrow \text{CHO} + \text{H}$ and $\text{CH}_2\text{O} + \text{h}\nu \rightarrow \text{CO} + \text{H}_2$ for Wavelengths From 240 to 360 nm

	03 COLUMN DENSITY					
	2.0E+12	5.0E+12	1.0E+13	2.0E+13	5.0E+13	1.0E+14
1.0E+13	2.58901E-04	2.58901E-04	2.58900E-04	2.58899E-04	2.58899E-04	2.58888E-04
2.0E+13	2.58901E-04	2.58901E-04	2.58900E-04	2.58899E-04	2.58899E-04	2.58888E-04
5.0E+13	2.58901E-04	2.58900E-04	2.58900E-04	2.58898E-04	2.58894E-04	2.58877E-04
1.0E+14	2.58900E-04	2.58900E-04	2.58899E-04	2.58898E-04	2.58894E-04	2.58873E-04
2.0E+14	2.58899E-04	2.58899E-04	2.58899E-04	2.58897E-04	2.58893E-04	2.58887E-04
5.0E+14	2.58897E-04	2.58896E-04	2.58896E-04	2.58894E-04	2.58890E-04	2.58883E-04
1.0E+15	2.58892E-04	2.58892E-04	2.58891E-04	2.58889E-04	2.58885E-04	2.58879E-04
2.0E+15	2.58883E-04	2.58882E-04	2.58882E-04	2.58880E-04	2.58876E-04	2.58872E-04
5.0E+15	2.58856E-04	2.58855E-04	2.58854E-04	2.58853E-04	2.58849E-04	2.58842E-04
1.0E+16	2.58812E-04	2.58811E-04	2.58811E-04	2.58809E-04	2.58805E-04	2.58798E-04
2.0E+16	2.58730E-04	2.58729E-04	2.58729E-04	2.58727E-04	2.58723E-04	2.58716E-04
5.0E+16	2.58523E-04	2.58522E-04	2.58522E-04	2.58520E-04	2.58516E-04	2.58509E-04
1.0E+17	2.58271E-04	2.58270E-04	2.58269E-04	2.58269E-04	2.58265E-04	2.58258E-04
2.0E+17	2.57962E-04	2.57961E-04	2.57961E-04	2.57959E-04	2.57955E-04	2.57949E-04
5.0E+17	2.57499E-04	2.57499E-04	2.57498E-04	2.57497E-04	2.57493E-04	2.57486E-04
1.0E+18	2.57029E-04	2.57028E-04	2.57028E-04	2.57026E-04	2.57022E-04	2.57016E-04
2.0E+18	2.56466E-04	2.56466E-04	2.56465E-04	2.56464E-04	2.56460E-04	2.56453E-04
5.0E+18	2.55942E-04	2.55941E-04	2.55941E-04	2.55939E-04	2.55935E-04	2.55929E-04
1.0E+19	2.55837E-04	2.55836E-04	2.55836E-04	2.55834E-04	2.55830E-04	2.55824E-04
2.0E+19	2.55824E-04	2.55824E-04	2.55823E-04	2.55822E-04	2.55818E-04	2.55811E-04
5.0E+19	2.55824E-04	2.55824E-04	2.55823E-04	2.55822E-04	2.55818E-04	2.55811E-04
1.0E+20	2.55824E-04	2.55824E-04	2.55823E-04	2.55822E-04	2.55818E-04	2.55811E-04
2.0E+20	2.55824E-04	2.55824E-04	2.55823E-04	2.55822E-04	2.55818E-04	2.55811E-04
5.0E+20	2.55824E-04	2.55824E-04	2.55823E-04	2.55822E-04	2.55818E-04	2.55811E-04
1.0E+21	2.55823E-04	2.55823E-04	2.55823E-04	2.55822E-04	2.55817E-04	2.55811E-04
2.0E+21	2.55818E-04	2.55818E-04	2.55818E-04	2.55817E-04	2.55812E-04	2.55805E-04
5.0E+21	2.55813E-04	2.55813E-04	2.55813E-04	2.55812E-04	2.55807E-04	2.55800E-04
1.0E+22	2.55823E-04	2.55823E-04	2.55822E-04	2.55821E-04	2.55817E-04	2.55811E-04
2.0E+22	2.55823E-04	2.55823E-04	2.55822E-04	2.55821E-04	2.55817E-04	2.55810E-04
5.0E+22	2.55821E-04	2.55821E-04	2.55820E-04	2.55820E-04	2.55819E-04	2.55815E-04
1.0E+23	2.55818E-04	2.55818E-04	2.55817E-04	2.55816E-04	2.55815E-04	2.55812E-04
2.0E+23	2.55813E-04	2.55813E-04	2.55813E-04	2.55812E-04	2.55811E-04	2.55808E-04
5.0E+23	2.55799E-04	2.55799E-04	2.55798E-04	2.55797E-04	2.55793E-04	2.55786E-04
1.0E+24	2.55781E-04	2.55781E-04	2.55780E-04	2.55780E-04	2.55779E-04	2.55775E-04
2.0E+24	2.55759E-04	2.55759E-04	2.55758E-04	2.55757E-04	2.55753E-04	2.55746E-04
5.0E+24	2.55739E-04	2.55739E-04	2.55738E-04	2.55737E-04	2.55733E-04	2.55727E-04
1.0E+25	2.55737E-04	2.55736E-04	2.55736E-04	2.55734E-04	2.55730E-04	2.55724E-04

02 COLUMN DENSITY

Table B18. Total Photodissociation Rate Coefficients for the Process $\text{CO}_2 + h\nu \rightarrow \text{CO} + \text{O}$ for Wavelengths From 166 to 215 nm (Contd)

		Q3 COLUMN DENSITY				
		5.0E+18	1.0E+19	2.0E+19	5.0E+19	1.0E+20
1.0E+13	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
2.0E+13	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
5.0E+13	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
1.0E+14	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
2.0E+14	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
5.0E+14	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
1.0E+15	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
2.0E+15	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
5.0E+15	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
1.0E+16	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
2.0E+16	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
5.0E+16	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
1.0E+17	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
2.0E+17	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
5.0E+17	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
1.0E+18	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
2.0E+18	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
5.0E+18	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
1.0E+19	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
2.0E+19	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
5.0E+19	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
1.0E+20	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
2.0E+20	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
5.0E+20	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
1.0E+21	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
2.0E+21	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
5.0E+21	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
1.0E+22	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
2.0E+22	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
5.0E+22	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
1.0E+23	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
2.0E+23	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
5.0E+23	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
1.0E+24	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40
2.0E+24	1.11969E-11	1.51540E-12	4.02055E-14	1.62200E-18	1.13333E-25	6.57859E-40

Table B18. Total Photodissociation Rate Coefficients for the Process $\text{CO}_2 + h\nu \rightarrow \text{CO} + \text{O}$ for Wavelengths From 166 to 215 nm (Contd)

	03 COLUMN DENSITY						02 COLUMN DEENSITY					
	2.0E+15	5.0E+15	1.0E+16	2.0E+16	5.0E+16	1.0E+17	2.0E+17	5.0E+17	1.0E+18	2.0E+18	5.0E+18	
1.0E+13	8.23559E-09	8.21569E-09	8.18264E-09	8.11694E-09	7.92300E-09	7.61006E-09	7.02095E-09	5.51462E-09	3.69031E-09	1.65855E-09		
2.0E+13	8.23553E-09	8.21563E-09	8.18258E-09	8.11688E-09	7.92294E-09	7.61000E-09	7.02090E-09	5.51458E-09	3.69028E-09	1.65854E-09		
5.0E+13	8.23535E-09	8.21546E-09	8.18241E-09	8.11671E-09	7.92277E-09	7.60984E-09	7.02075E-09	5.51446E-09	3.69021E-09	1.65850E-09		
1.0E+14	8.23506E-09	8.21516E-09	8.18211E-09	8.11642E-09	7.92249E-09	7.60957E-09	7.02050E-09	5.51427E-09	3.69008E-09	1.65845E-09		
2.0E+14	8.23447E-09	8.21457E-09	8.18153E-09	8.11584E-09	7.92192E-09	7.60902E-09	7.02000E-09	5.51388E-09	3.68982E-09	1.65833E-09		
5.0E+14	8.23270E-09	8.21281E-09	8.17977E-09	8.11409E-09	7.92022E-09	7.60739E-09	7.01849E-09	5.51270E-09	3.68930E-09	1.65793E-09		
1.0E+15	8.22975E-09	8.20987E-09	8.17784E-09	8.11119E-09	7.91739E-09	7.60467E-09	7.01599E-09	5.51074E-09	3.68773E-09	1.65741E-09		
2.0E+15	8.22386E-09	8.20400E-09	8.17099E-09	8.10538E-09	7.91172E-09	7.59923E-09	7.01098E-09	5.50682E-09	3.68513E-09	1.65626E-09		
5.0E+15	8.20623E-09	8.18641E-09	8.15347E-09	8.08801E-09	7.89477E-09	7.58296E-09	6.99598E-09	5.49508E-09	3.67733E-09	1.65282E-09		
1.0E+16	8.17696E-09	8.15721E-09	8.12440E-09	8.05917E-09	7.866633E-09	7.55595E-09	6.97109E-09	5.47561E-09	3.66439E-09	1.64711E-09		
2.0E+16	8.11889E-09	8.09928E-09	8.06670E-09	8.0194E-09	7.81070E-09	7.50234E-09	6.92170E-09	5.43696E-09	3.63871E-09	1.63577E-09		
5.0E+16	7.94821E-09	7.92902E-09	7.89714E-09	7.83376E-09	7.64668E-09	7.60709E-09	6.94811E-09	5.32339E-09	3.56326E-09	1.60246E-09		
1.0E+17	7.67522E-09	7.65670E-09	7.62593E-09	7.56476E-09	7.38421E-09	7.09286E-09	6.54437E-09	5.14173E-09	3.44258E-09	1.54920E-09		
2.0E+17	7.15081E-09	7.13357E-09	7.10493E-09	7.04801E-09	6.87997E-09	6.6081E-09	6.09833E-09	4.79270E-09	3.21064E-09	1.44677E-09		
5.0E+17	5.85974E-09	5.84567E-09	5.82229E-09	5.77581E-09	5.63981E-09	5.41720E-09	5.00309E-09	3.93360E-09	2.63989E-09	1.19483E-09		
1.0E+18	4.38135E-09	4.37090E-09	4.35155E-09	4.31105E-09	4.21719E-09	4.05279E-09	3.74314E-09	2.95019E-09	2.98677E-09	9.06714E-10		
2.0E+18	2.83407E-09	2.82743E-09	2.81639E-09	2.79445E-09	2.72966E-09	2.62504E-09	2.42783E-09	1.92180E-09	1.30434E-09	6.06124E-10		
5.0E+18	1.55073E-09	1.54725E-09	1.54147E-09	1.52998E-09	1.49604E-09	1.44116E-09	1.33751E-09	1.92180E-09	1.39870E-10	3.58122E-10		
1.0E+19	1.14257E-09	1.14010E-09	1.13598E-09	1.12780E-09	1.10361E-09	1.06448E-09	9.90439E-10	7.98532E-10	5.59504E-10	2.78444E-10		
2.0E+19	9.20991E-10	9.19045E-10	9.15810E-10	9.09377E-10	8.90355E-10	8.59560E-10	8.01226E-10	6.49564E-10	4.59465E-10	2.33258E-10		
5.0E+19	7.11624E-10	7.10164E-10	7.07738E-10	7.02911E-10	6.88636E-10	6.65510E-10	6.21645E-10	5.07214E-10	3.62768E-10	1.88543E-10		
1.0E+20	5.73567E-10	5.72422E-10	5.66731E-10	5.55254E-10	5.37359E-10	5.02866E-10	4.12617E-10	2.97993E-10	1.58086E-10	6.47265E-10		
2.0E+20	4.47148E-10	4.46286E-10	4.44854E-10	4.42003E-10	4.33565E-10	4.19877E-10	3.93849E-10	3.25484E-10	2.37963E-10	1.29490E-10		
5.0E+20	4.46927E-10	4.46066E-10	4.44634E-10	4.41784E-10	4.19669E-10	3.93652E-10	3.25319E-10	2.52319E-10	2.37839E-10	1.29494E-10		
1.0E+21	3.08433E-10	3.07875E-10	3.05101E-10	3.06947E-10	2.99633E-10	2.90751E-10	2.73817E-10	2.29036E-10	1.70898E-10	9.68980E-11		
2.0E+21	2.34369E-10	2.33997E-10	2.31968E-10	2.28025E-10	2.21612E-10	2.09361E-10	1.76783E-10	1.34005E-10	1.78375E-11	7.8375E-11		
5.0E+21	1.84499E-10	1.84196E-10	1.83691E-10	1.82687E-10	1.79710E-10	1.74864E-10	1.65593E-10	1.40837E-10	1.08055E-10	6.47265E-11		
1.0E+22	1.41710E-10	1.41484E-10	1.41109E-10	1.40362E-10	1.38148E-10	1.34542E-10	1.27637E-10	1.09152E-10	8.45402E-11	5.16358E-11		
2.0E+22	1.15859E-10	1.15677E-10	1.15373E-10	1.14768E-10	1.12973E-10	1.10051E-10	1.04455E-10	8.94741E-11	6.95121E-11	4.27606E-11		
5.0E+22	8.48856E-11	8.47527E-11	8.45161E-11	8.40914E-11	8.278862E-11	8.06608E-11	7.65907E-11	6.56940E-11	5.11706E-11	3.16775E-11		
1.0E+23	4.98125E-11	4.97335E-11	4.96020E-11	4.93402E-11	4.85642E-11	4.73012E-11	4.48B46E-11	3.84294E-11	2.98619E-11	1.84374E-11		
2.0E+23	2.07956E-11	2.07625E-11	2.07075E-11	2.05980E-11	2.02734E-11	1.97452E-11	1.87351E-11	1.60400E-11	1.24702E-11	1.72202E-12		
5.0E+23	3.00891E-12	3.00448E-12	3.00111E-12	2.98243E-12	2.93890E-12	2.86802E-12	2.73219E-12	2.36794E-12	1.88005E-12	1.2156E-12		
1.0E+24	4.35417E-13	4.34426E-13	4.32780E-13	4.29508E-13	4.19844E-13	4.04232E-13	3.74777E-13	2.99008E-13	2.06060E-13	9.9738E-14		
2.0E+24	3.94567E-13	3.93629E-13	3.92071E-13	3.88973E-13	3.79825E-13	3.65056E-13	3.47227E-13	2.65877E-13	1.79013E-13	8.13910E-14		

Table B18. Total Photodissociation Rate Coefficients for the Process $\text{CO}_2 + \text{h}\nu \rightarrow \text{CO} + \text{O}$ for Wavelengths From 166 to 215 nm

	03 COLUMN DENSITY									
	1.0E+12	2.0E+12	5.0E+12	1.0E+13	2.0E+13	5.0E+13	1.0E+14	2.0E+14	5.0E+14	1.0E+15
1.0E+13	8.24887E-09	8.24886E-09	8.24884E-09	8.24881E-09	8.24875E-09	8.24875E-09	8.24875E-09	8.24875E-09	8.24875E-09	8.24875E-09
2.0E+13	8.24881E-09	8.24881E-09	8.24879E-09	8.24879E-09	8.24875E-09	8.24875E-09	8.24875E-09	8.24875E-09	8.24875E-09	8.24875E-09
5.0E+13	8.24864E-09	6.24863E-09	8.24861E-09	8.24861E-09	8.24858E-09	8.24851E-09	8.24831E-09	8.24798E-09	8.24731E-09	8.24532E-09
1.0E+14	8.24834E-09	8.24833E-09	8.24832E-09	8.24831E-09	8.24828E-09	8.24821E-09	8.24801E-09	8.24768E-09	8.24702E-09	8.24702E-09
2.0E+14	8.24775E-09	6.24774E-09	8.24774E-09	8.24774E-09	8.24769E-09	8.24762E-09	8.24747E-09	8.24709E-09	8.24643E-09	8.24443E-09
5.0E+14	8.24598E-09	8.24597E-09	8.24595E-09	8.24595E-09	8.24592E-09	8.24585E-09	8.24565E-09	8.24532E-09	8.24266E-09	8.24111E-09
1.0E+15	8.24363E-09	8.24302E-09	8.24300E-09	8.24297E-09	8.24290E-09	8.24290E-09	8.24270E-09	8.24237E-09	8.24170E-09	8.23971E-09
2.0E+15	8.23713E-09	8.23712E-09	8.23710E-09	8.23707E-09	8.23700E-09	8.23660E-09	8.23647E-09	8.23580E-09	8.23381E-09	8.23049E-09
5.0E+15	8.21947E-09	8.21946E-09	8.21945E-09	8.21944E-09	8.21943E-09	8.21943E-09	8.21943E-09	8.21881E-09	8.21616E-09	8.21285E-09
1.0E+16	8.19015E-09	8.19015E-09	8.19013E-09	8.19013E-09	8.19003E-09	8.19003E-09	8.19003E-09	8.18983E-09	8.18844E-09	8.18356E-09
2.0E+16	8.11985E-09	8.11985E-09	8.11985E-09	8.11985E-09	8.11985E-09	8.11985E-09	8.11985E-09	8.13135E-09	8.13067E-09	8.12543E-09
5.0E+16	7.96102E-09	7.96102E-09	7.96100E-09	7.96097E-09	7.96090E-09	7.96090E-09	7.96071E-09	7.96039E-09	7.95975E-09	7.95782E-09
1.0E+17	7.68759E-09	7.68758E-09	7.68756E-09	7.68753E-09	7.68747E-09	7.68729E-09	7.68698E-09	7.68636E-09	7.68450E-09	7.68141E-09
2.0E+17	7.16231E-09	7.16229E-09	7.16226E-09	7.16223E-09	7.16220E-09	7.16203E-09	7.16174E-09	7.16117E-09	7.15944E-09	7.15656E-09
5.0E+17	5.86913E-09	5.86913E-09	5.86912E-09	5.86912E-09	5.86905E-09	5.86905E-09	5.86890E-09	5.86867E-09	5.86679E-09	5.86444E-09
1.0E+18	4.38832E-09	4.38832E-09	4.38831E-09	4.38831E-09	4.38826E-09	4.38826E-09	4.38815E-09	4.38798E-09	4.38763E-09	4.38658E-09
2.0E+18	2.83851E-09	2.83850E-09	2.83849E-09	2.83849E-09	2.83840E-09	2.83840E-09	2.83840E-09	2.83807E-09	2.83740E-09	2.83629E-09
5.0E+18	1.55305E-09	1.55304E-09	1.55304E-09	1.55304E-09	1.55299E-09	1.55299E-09	1.55299E-09	1.55293E-09	1.55282E-09	1.55189E-09
1.0E+19	1.14423E-09	1.14423E-09	1.14422E-09	1.14422E-09	1.14421E-09	1.14421E-09	1.14421E-09	1.14419E-09	1.14415E-09	1.14340E-09
2.0E+19	9.22598E-10	9.22598E-10	9.22264E-10	9.22264E-10	9.22278E-10	9.22258E-10	9.22258E-10	9.22160E-10	9.21965E-10	9.21640E-10
5.0E+19	7.12598E-10	7.12598E-10	7.12597E-10	7.12597E-10	7.12594E-10	7.12574E-10	7.12550E-10	7.12535E-10	7.12111E-10	7.11211E-10
1.0E+20	5.74331E-10	5.74331E-10	5.74330E-10	5.74330E-10	5.74328E-10	5.74328E-10	5.74313E-10	5.74256E-10	5.74141E-10	5.73950E-10
2.0E+20	4.47723E-10	4.47723E-10	4.47722E-10	4.47722E-10	4.47718E-10	4.47718E-10	4.47709E-10	4.47695E-10	4.47662E-10	4.47436E-10
5.0E+20	4.47502E-10	4.47502E-10	4.47501E-10	4.47501E-10	4.47497E-10	4.47497E-10	4.47498E-10	4.47445E-10	4.47445E-10	4.47215E-10
1.0E+21	3.08805E-10	3.08805E-10	3.06804E-10	3.06804E-10	3.08803E-10	3.08803E-10	3.08802E-10	3.08796E-10	3.08787E-10	3.08619E-10
2.0E+21	2.34638E-10	2.34638E-10	2.34637E-10	2.34637E-10	2.34636E-10	2.34636E-10	2.34631E-10	2.34624E-10	2.34571E-10	2.34504E-10
5.0E+21	1.84702E-10	1.84702E-10	1.84701E-10	1.84701E-10	1.84701E-10	1.84701E-10	1.84693E-10	1.84692E-10	1.84651E-10	1.84600E-10
1.0E+22	1.41865E-10	1.41865E-10	1.41860E-10	1.41860E-10	1.41859E-10	1.41857E-10	1.41857E-10	1.41853E-10	1.41853E-10	1.41795E-10
2.0E+22	1.15981E-10	1.15981E-10	1.15981E-10	1.15981E-10	1.15980E-10	1.15978E-10	1.15975E-10	1.15969E-10	1.15951E-10	1.15920E-10
5.0E+22	8.49743E-11	8.49743E-11	8.49743E-11	8.49743E-11	8.49739E-11	8.49735E-11	8.49721E-11	8.49699E-11	8.49655E-11	8.49300E-11
1.0E+23	4.98653E-11	4.98653E-11	4.98652E-11	4.98652E-11	4.98651E-11	4.98648E-11	4.98640E-11	4.98627E-11	4.98521E-11	4.98389E-11
2.0E+23	2.08177E-11	2.08177E-11	2.08176E-11	2.08176E-11	2.08176E-11	2.08174E-11	2.08171E-11	2.08155E-11	2.08145E-11	2.08066E-11
5.0E+23	3.01187E-12	3.01186E-12	3.01186E-12	3.01186E-12	3.01185E-12	3.01179E-12	3.01175E-12	3.01172E-12	3.01133E-12	3.01102E-12
1.0E+24	4.36077E-13	4.36078E-13	4.36075E-13	4.36075E-13	4.36062E-13	4.36061E-13	4.36045E-13	4.35747E-13	4.35591E-13	4.35374E-13
2.0E+24	3.95193E-13	3.95192E-13	3.95193E-13	3.95192E-13	3.95190E-13	3.95187E-13	3.95183E-13	3.95178E-13	3.95173E-13	3.95037E-13

Table B17. Total Photodissociation Rate Coefficients for the Process $\text{CO}_2 + h\nu \rightarrow \text{CO} + \text{O}^1\text{D}$ for Wavelengths From 116.2 to 166 nm (Contd)

03 COLUMN DENSITY						
	5.0E+17	1.0E+18	2.0E+18	5.0E+18	1.0E+19	2.0E+19
1.0E+13	1.44282E-08	6.24892E-09	1.78563E-09	8.73281E-11	8.97046E-13	1.29487E-16
2.0E+13	1.44273E-08	6.24861E-09	1.78556E-09	8.73257E-11	8.97025E-13	1.29484E-16
5.0E+13	1.44245E-08	6.24766E-09	1.78536E-09	8.73182E-11	8.96959E-13	1.29476E-16
1.0E+14	1.44198E-08	6.24608E-09	1.78502E-09	8.73058E-11	8.96849E-13	1.29461E-16
2.0E+14	1.44104E-06	6.24293E-09	1.78433E-09	8.72810E-11	8.96630E-13	1.29433E-16
5.0E+14	1.43823E-08	6.23348E-09	1.78228E-09	8.72066E-11	8.95973E-13	1.29348E-16
1.0E+15	1.43356E-08	6.21777E-09	1.77887E-09	8.70829E-11	8.94879E-13	1.29205E-16
2.0E+15	1.42429E-08	6.18650E-09	1.77208E-09	8.68360E-11	8.92695E-13	1.28921E-16
5.0E+15	1.39692E-08	6.09393E-09	1.75189E-09	8.60999E-11	8.86177E-13	1.28073E-16
1.0E+16	1.35281E-08	5.94358E-09	1.71889E-09	8.48891E-11	8.75429E-13	1.26673E-16
2.0E+16	1.26987E-08	5.65705E-09	1.65517E-09	8.25252E-11	8.54351E-13	1.23919E-16
5.0E+16	1.05782E-08	4.89851E-09	1.48097E-09	7.58722E-11	7.94332E-13	1.16022E-16
1.0E+17	7.97646E-09	3.90476E-09	1.23854E-09	6.60946E-11	7.04143E-13	1.03993E-16
2.0E+17	4.87398E-09	2.58936E-09	8.85155E-10	5.05224E-11	5.55023E-13	8.36308E-17
5.0E+17	1.62428E-09	9.52031E-10	3.65897E-10	2.36228E-11	2.77402E-13	4.37894E-17
1.0E+18	4.54105E-10	2.58205E-10	1.04641E-10	7.34631E-12	9.15228E-14	1.51281E-17
2.0E+18	8.08768E-11	2.89490E-11	1.13689E-11	8.34787E-13	1.03426E-14	1.85701E-18
5.0E+18	1.89959E-11	1.01080E-12	2.41356E-14	1.57727E-15	2.07641E-17	3.59944E-21
1.0E+19	8.09561E-12	4.35639E-13	1.41722E-15	4.80283E-20	6.31712E-22	1.09529E-25
2.0E+19	2.38741E-12	1.22251E-13	4.06649E-16	1.63975E-23	5.85059E-31	1.01439E-34
5.0E+19	2.84047E-13	5.74015E-15	1.43921E-17	5.76535E-25	2.94718E-37	1.57837E-61
1.0E+20	1.07654E-13	6.66417E-16	3.14633E-19	2.36667E-27	1.20319E-39	3.15692E-64
2.0E+20	3.14917E-14	8.62465E-17	3.49187E-20	2.43571E-30	2.01228E-44	5.27258E-69
5.0E+20	1.36298E-15	2.23290E-19	8.63655E-23	5.94019E-33	6.85986E-50	3.37128E-83
1.0E+21	9.11393E-18	9.31033E-23	3.92055E-27	2.69684E-37	3.11437E-54	4.15337E-88
2.0E+21	4.13750E-22	3.79249E-27	8.39948E-36	5.55861E-46	6.41920E-63	8.56074E-97
5.0E+21	3.87172E-35	3.54885E-40	2.98165E-50	4.86740E-72	5.62098E-89	7.49622E-123
1.0E+22	7.46758E-57	6.84485E-62	5.75085E-72	3.41064E-102	2.09105E-132	2.78865E-268
2.0E+22	2.77800E-100	2.54634E-105	2.13936E-115	1.26687E-145	5.31163E-196	3.85921E-253
5.0E+22	1.43017E-230	1.31090E-235	1.10139E-245	6.53196E-276	0.	0.

02 COLUMN DENSITY

Table B17. Total Photodissociation Rate Coefficients for the Process $\text{CO}_2 + h\nu \rightarrow \text{CO} + \text{O}(\text{D})$ for Wavelengths From 116.2 to 166 nm (Contd)

03 COLUMN DENSITY						
2.0E+14	5.0E+14	1.0E+15	2.0E+15	5.0E+15	1.0E+16	2.0E+16
1.0E+13	8.82221E-08	8.79581E-08	8.75211E-08	8.66585E-08	8.41591E-08	8.02701E-08
2.0E+13	8.82167E-08	8.79527E-08	8.75157E-08	8.66532E-08	8.41539E-08	8.02650E-08
5.0E+13	8.82006E-08	8.79366E-08	8.74997E-08	8.66372E-08	8.41381E-08	8.02496E-08
1.0E+14	8.81738E-08	8.79098E-08	8.74730E-08	8.66107E-08	8.41119E-08	8.02240E-08
2.0E+14	8.81202E-08	8.78563E-08	8.74196E-08	8.65575E-08	8.40596E-08	8.01729E-08
5.0E+14	8.79597E-08	8.76960E-08	8.72597E-08	8.63984E-08	8.39028E-08	8.00200E-08
1.0E+15	8.76933E-08	8.74301E-08	8.69944E-08	8.61344E-08	8.36427E-08	8.79761E-08
2.0E+15	8.71649E-08	8.69024E-08	8.64681E-08	8.56107E-08	8.31267E-08	8.92625E-08
5.0E+15	8.56136E-08	8.53534E-08	8.49229E-08	8.40732E-08	8.16116E-08	7.77840E-08
1.0E+16	8.31372E-08	8.28807E-08	8.24563E-08	8.16187E-08	8.91930E-08	7.54235E-08
2.0E+16	7.85653E-08	7.83156E-08	7.79024E-08	7.70871E-08	7.47276E-08	7.10652E-08
5.0E+16	6.74127E-08	6.67935E-08	6.547151E-08	6.43595E-08	6.38334E-08	6.04304E-08
1.0E+17	5.49301E-08	5.43595E-08	5.36587E-08	5.16375E-08	5.07761E-08	4.85219E-08
2.0E+17	4.21191E-08	4.19228E-08	4.15984E-08	4.09596E-08	3.91210E-08	3.62994E-08
5.0E+17	3.12219E-08	3.10436E-08	3.07491E-08	3.01694E-08	2.85041E-08	2.59585E-08
1.0E+18	2.69793E-08	2.68113E-08	2.65338E-08	2.59879E-08	2.44211E-08	2.20304E-08
2.0E+18	2.42727E-08	2.41140E-08	2.38519E-08	2.33365E-08	2.18582E-08	1.96063E-08
5.0E+18	2.16428E-08	2.14958E-08	2.12531E-08	2.07760E-08	1.94088E-08	1.73297E-08
1.0E+19	1.99580E-08	1.98205E-08	1.95934E-08	1.91471E-08	1.78686E-08	1.59254E-08
2.0E+19	1.78972E-08	1.77733E-08	1.75687E-08	1.71666E-08	1.60146E-08	1.42642E-08
5.0E+19	1.32281E-08	1.31364E-08	1.29850E-08	1.26873E-08	1.18346E-08	1.05338E-08
1.0E+20	8.02109E-09	7.96546E-09	7.87361E-09	7.69307E-09	7.17592E-09	6.39010E-09
2.0E+20	2.95034E-09	2.92988E-09	2.89609E-09	2.82968E-09	2.63944E-09	2.35037E-09
5.0E+20	1.46878E-10	1.45860E-10	1.44178E-10	1.40871E-10	1.31400E-10	1.17008E-10
1.0E+21	9.89656E-13	9.82792E-13	9.71457E-13	9.49179E-13	8.85363E-13	7.88393E-13
2.0E+21	4.49303E-17	4.46187E-17	4.41041E-17	4.30927E-17	4.01954E-17	3.57930E-17
5.0E+21	4.20441E-30	4.17525E-30	4.12709E-30	4.03245E-30	3.76133E-30	3.34937E-30
1.0E+22	8.10925E-52	8.05301E-52	7.96013E-52	7.77758E-52	7.25467E-52	6.46011E-52
2.0E+22	3.01670E-95	2.99578E-95	2.96123E-95	2.89332E-95	2.69879E-95	2.40321E-95
5.0E+22	1.55306E-225	1.54229E-225	1.52450E-225	1.48954E-225	1.38939E-225	1.23722E-225

02 COLUMN DENSITY						
2.0E+14	5.0E+14	1.0E+15	2.0E+15	5.0E+15	1.0E+16	2.0E+16
1.0E+13	8.4055E-08	8.34210E-08	8.26550E-08	8.13456E-08	8.02650E-08	7.34056E-08
2.0E+13	8.37669E-08	8.31210E-08	8.23082E-08	8.13391E-08	8.02496E-08	7.33910E-08
5.0E+13	8.37567E-C8	8.31210E-08	8.23082E-08	8.13391E-08	8.02496E-08	7.337395E-C8
1.0E+14	8.37809E-08	8.31210E-08	8.23082E-08	8.13366E-08	8.02240E-08	7.33179E-08
2.0E+14	8.37853E-08	8.31210E-08	8.23082E-08	8.13179E-08	8.01729E-08	7.33179E-08
5.0E+14	8.37853E-08	8.31210E-08	8.23082E-08	8.13179E-08	8.01729E-08	7.33179E-08
1.0E+15	8.37853E-08	8.31210E-08	8.23082E-08	8.13179E-08	8.01729E-08	7.33179E-08
2.0E+15	8.37853E-08	8.31210E-08	8.23082E-08	8.13179E-08	8.01729E-08	7.33179E-08
5.0E+15	8.37853E-08	8.31210E-08	8.23082E-08	8.13179E-08	8.01729E-08	7.33179E-08
1.0E+16	8.37853E-08	8.31210E-08	8.23082E-08	8.13179E-08	8.01729E-08	7.33179E-08
2.0E+16	8.37853E-08	8.31210E-08	8.23082E-08	8.13179E-08	8.01729E-08	7.33179E-08
5.0E+16	8.37853E-08	8.31210E-08	8.23082E-08	8.13179E-08	8.01729E-08	7.33179E-08
1.0E+17	8.37853E-08	8.31210E-08	8.23082E-08	8.13179E-08	8.01729E-08	7.33179E-08
2.0E+17	8.37853E-08	8.31210E-08	8.23082E-08	8.13179E-08	8.01729E-08	7.33179E-08
5.0E+17	8.37853E-08	8.31210E-08	8.23082E-08	8.13179E-08	8.01729E-08	7.33179E-08
1.0E+18	8.37853E-08	8.31210E-08	8.23082E-08	8.13179E-08	8.01729E-08	7.33179E-08
2.0E+18	8.37853E-08	8.31210E-08	8.23082E-08	8.13179E-08	8.01729E-08	7.33179E-08
5.0E+18	8.37853E-08	8.31210E-08	8.23082E-08	8.13179E-08	8.01729E-08	7.33179E-08
1.0E+19	8.37853E-08	8.31210E-08	8.23082E-08	8.13179E-08	8.01729E-08	7.33179E-08
2.0E+19	8.37853E-08	8.31210E-08	8.23082E-08	8.13179E-08	8.01729E-08	7.33179E-08
5.0E+19	8.37853E-08	8.31210E-08	8.23082E-08	8.13179E-08	8.01729E-08	7.33179E-08
1.0E+20	8.37853E-08	8.31210E-08	8.23082E-08	8.13179E-08	8.01729E-08	7.33179E-08
2.0E+20	8.37853E-08	8.31210E-08	8.23082E-08	8.13179E-08	8.01729E-08	7.33179E-08
5.0E+20	8.37853E-08	8.31210E-08	8.23082E-08	8.13179E-08	8.01729E-08	7.33179E-08
1.0E+21	8.37853E-08	8.31210E-08	8.23082E-08	8.13179E-08	8.01729E-08	7.33179E-08
2.0E+21	8.37853E-08	8.31210E-08	8.23082E-08	8.13179E-08	8.01729E-08	7.33179E-08
5.0E+21	8.37853E-08	8.31210E-08	8.23082E-08	8.13179E-08	8.01729E-08	7.33179E-08
1.0E+22	8.37853E-08	8.31210E-08	8.23082E-08	8.13179E-08	8.01729E-08	7.33179E-08
2.0E+22	8.37853E-08	8.31210E-08	8.23082E-08	8.13179E-08	8.01729E-08	7.33179E-08
5.0E+22	8.37853E-08	8.31210E-08	8.23082E-08	8.13179E-08	8.01729E-08	7.33179E-08

Table B17. Total Photodissociation Rate Coefficients for the Process $\text{CO}_2 + h\nu \rightarrow \text{CO} + \text{O}^1\text{D}$ for Wavelengths From 116.2 to 166 nm

	O3 COLUMN DENSITY									
	1.0E+11	2.0E+11	5.0E+11	1.0E+12	2.0E+12	5.0E+12	1.0E+13	2.0E+13	5.0E+13	1.0E+14
1.0E+13	8.83988E-08	8.83987E-08	8.83984E-08	8.83980E-08	8.83971E-08	8.83945E-08	8.83900E-08	8.83812E-08	8.83546E-08	8.83104E-08
2.0E+13	8.83934E-08	8.83933E-08	8.83931E-08	8.83926E-08	8.83917E-08	8.83891E-08	8.83847E-08	8.83758E-08	8.83493E-08	8.83050E-08
5.0E+13	8.83773E-08	8.83772E-08	8.83769E-08	8.83765E-08	8.83760E-08	8.83685E-08	8.83597E-08	8.83331E-08	8.82889E-08	8.82261E-08
1.0E+14	8.83504E-08	8.83504E-08	8.83501E-08	8.83496E-08	8.83491E-08	8.83446E-08	8.83417E-08	8.83328E-08	8.83063E-08	8.82621E-08
2.0E+14	8.82968E-08	8.82967E-08	8.82964E-08	8.82960E-08	8.82951E-08	8.82924E-08	8.82912E-08	8.82526E-08	8.82084E-08	8.82084E-08
5.0E+14	8.81361E-08	8.81360E-08	8.81357E-08	8.81353E-08	8.81344E-08	8.81318E-08	8.81273E-08	8.81185E-08	8.80920E-08	8.80478E-08
1.0E+15	8.78695E-08	8.78694E-08	8.78691E-08	8.78687E-08	8.78684E-08	8.78678E-08	8.78671E-08	8.78519E-08	8.78254E-08	8.77814E-08
2.0E+15	8.73405E-08	8.73405E-08	8.73402E-08	8.73398E-08	8.73395E-08	8.73392E-08	8.73318E-08	8.73230E-08	8.72966E-08	8.72527E-08
5.0E+15	8.57877E-08	8.57876E-08	8.57874E-08	8.57869E-08	8.57861E-08	8.57834E-08	8.57791E-08	8.57703E-08	8.57442E-08	8.57006E-08
1.0E+16	8.33088E-08	8.33087E-08	8.33085E-08	8.33083E-08	8.33081E-08	8.33072E-08	8.33046E-08	8.33003E-08	8.32917E-08	8.32659E-08
2.0E+16	7.87324E-08	7.87323E-08	7.87321E-08	7.87319E-08	7.87317E-08	7.87308E-08	7.87283E-08	7.87241E-08	7.87157E-08	7.86906E-08
5.0E+16	6.75689E-08	6.75688E-08	6.75687E-08	6.75686E-08	6.75685E-08	6.75674E-08	6.75651E-08	6.75611E-08	6.75298E-08	6.74908E-08
1.0E+17	5.50741E-08	5.50741E-08	5.50739E-08	5.50739E-08	5.50735E-08	5.50728E-08	5.50705E-08	5.50670E-08	5.50598E-08	5.50381E-08
2.0E+17	4.22505E-08	4.22505E-08	4.22503E-08	4.22499E-08	4.22493E-08	4.22473E-08	4.22440E-08	4.22374E-08	4.22177E-08	4.21848E-08
5.0E+17	3.13414E-08	3.13413E-08	3.13412E-08	3.13409E-08	3.13403E-08	3.13385E-08	3.13355E-08	3.13303E-08	3.132917E-08	3.12816E-08
1.0E+18	2.70919E-08	2.70919E-08	2.70917E-08	2.70914E-08	2.70912E-08	2.70908E-08	2.70891E-08	2.70863E-08	2.70807E-08	2.70356E-08
2.0E+18	2.43791E-08	2.43790E-08	2.43789E-08	2.43786E-08	2.43781E-08	2.43766E-08	2.43738E-08	2.43685E-08	2.43525E-08	2.43259E-08
5.0E+18	2.17413E-08	2.17412E-08	2.17411E-08	2.17408E-08	2.17403E-08	2.17388E-08	2.17364E-08	2.17314E-08	2.17166E-08	2.16921E-08
1.0E+19	2.00502E-08	2.00501E-08	2.00500E-08	2.00497E-08	2.00493E-08	2.00479E-08	2.00456E-08	2.00410E-08	2.00271E-08	2.00041E-08
2.0E+19	1.79802E-08	1.79802E-08	1.79801E-08	1.79798E-08	1.79794E-08	1.79782E-08	1.79761E-08	1.79719E-08	1.79595E-08	1.79387E-08
5.0E+19	1.32896E-08	1.32896E-08	1.32895E-08	1.32893E-08	1.32891E-08	1.32890E-08	1.32881E-08	1.32865E-08	1.32835E-08	1.32742E-08
1.0E+20	8.05837E-09	8.05835E-09	8.05829E-09	8.05820E-09	8.05801E-09	8.05745E-09	8.05652E-09	8.04905E-09	8.03972E-09	8.03972E-09
2.0E+20	2.96405E-09	2.96405E-09	2.96403E-09	2.96400E-09	2.96393E-09	2.96372E-09	2.96338E-09	2.96269E-09	2.96063E-09	2.95720E-09
5.0E+20	1.47561E-10	1.47561E-10	1.47560E-10	1.47558E-10	1.47555E-10	1.47552E-10	1.47544E-10	1.47493E-10	1.47390E-10	1.47220E-10
1.0E+21	9.94256E-13	9.94254E-13	9.94247E-13	9.94235E-13	9.94212E-13	9.94143E-13	9.94028E-13	9.93797E-13	9.93106E-13	9.91954E-13
2.0E+21	4.51392E-17	4.51391E-17	4.51387E-17	4.51382E-17	4.51372E-17	4.51340E-17	4.51183E-17	4.50869E-17	4.50347E-17	
5.0E+21	4.22395E-30	4.22394E-30	4.22391E-30	4.22386E-30	4.22377E-30	4.22347E-30	4.22200E-30	4.21907E-30	4.21417E-30	
1.0E+22	8.14695E-52	8.14693E-52	8.14687E-52	8.14678E-52	8.14659E-52	8.14602E-52	8.14319E-52	8.12809E-52	8.12752E-52	
2.0E+22	3.03073E-95	3.03072E-95	3.03070E-95	3.03066E-95	3.03059E-95	3.03038E-95	3.03003E-95	3.02722E-95	3.02371E-95	
5.0E+22	1.56028E-225	1.56027E-225	1.56026E-225	1.56025E-225	1.56021E-225	1.56010E-225	1.55992E-225	1.55956E-225	1.55847E-225	1.55667E-225

Table B16. Total Photodissociation Rate Coefficients for the Process $\text{CH}_4 + h\nu \rightarrow \text{CH}_3 + \text{H}$ and $\text{CH}_4 + h\nu \rightarrow \text{CH}_2 + \text{H}_2$ for Wavelengths From 116 to 164 nm (Contd)

	02 COLUMN DENSITY		03 COLUMN DENSITY	
	$5.0\text{E}+16$	$1.0\text{E}+17$	$2.0\text{E}+17$	$5.0\text{E}+17$
$1.0\text{E}+13$	$2.97405\text{E}-06$	$1.00972\text{E}-06$	$1.47022\text{E}-07$	$6.68945\text{E}-09$
$2.0\text{E}+13$	$2.97404\text{E}-06$	$1.00971\text{E}-06$	$1.47020\text{E}-07$	$6.68929\text{E}-09$
$5.0\text{E}+13$	$2.97402\text{E}-06$	$1.00970\text{E}-06$	$1.47014\text{E}-07$	$6.68880\text{E}-09$
$1.0\text{E}+14$	$2.97397\text{E}-06$	$1.00967\text{E}-06$	$1.47005\text{E}-07$	$6.68799\text{E}-09$
$2.0\text{E}+14$	$2.97388\text{E}-06$	$1.00962\text{E}-06$	$1.46986\text{E}-07$	$6.68636\text{E}-09$
$5.0\text{E}+14$	$2.97362\text{E}-06$	$1.00947\text{E}-06$	$1.46929\text{E}-07$	$6.68149\text{E}-09$
$1.0\text{E}+15$	$2.97313\text{E}-06$	$1.00922\text{E}-06$	$1.46836\text{E}-07$	$6.67341\text{E}-09$
$2.0\text{E}+15$	$2.97231\text{E}-06$	$1.00872\text{E}-06$	$1.46650\text{E}-07$	$6.65742\text{E}-09$
$5.0\text{E}+15$	$2.96975\text{E}-06$	$1.00726\text{E}-06$	$1.46106\text{E}-07$	$6.61059\text{E}-09$
$1.0\text{E}+16$	$2.96556\text{E}-06$	$1.00493\text{E}-06$	$1.45243\text{E}-07$	$6.53631\text{E}-09$
$2.0\text{E}+16$	$2.95803\text{E}-06$	$1.00064\text{E}-06$	$1.43663\text{E}-07$	$6.40065\text{E}-09$
$5.0\text{E}+16$	$2.93901\text{E}-06$	$9.90083\text{E}-07$	$1.39868\text{E}-07$	$6.07711\text{E}-09$
$1.0\text{E}+17$	$2.91601\text{E}-06$	$9.77741\text{E}-07$	$1.35633\text{E}-07$	$5.72204\text{E}-09$
$2.0\text{E}+17$	$2.88619\text{E}-06$	$9.62535\text{E}-07$	$1.30789\text{E}-07$	$5.32920\text{E}-09$
$5.0\text{E}+17$	$2.83733\text{E}-06$	$9.36986\text{E}-07$	$1.23134\text{E}-07$	$4.71516\text{E}-09$
$1.0\text{E}+18$	$2.77549\text{E}-06$	$9.09088\text{E}-07$	$1.14694\text{E}-07$	$3.99244\text{E}-09$
$2.0\text{E}+18$	$2.69502\text{E}-06$	$8.71894\text{E}-07$	$1.03686\text{E}-07$	$2.99423\text{E}-09$
$5.0\text{E}+18$	$2.55052\text{E}-06$	$8.11447\text{E}-07$	$8.10388\text{E}-08$	$1.56412\text{E}-09$
$1.0\text{E}+19$	$2.39877\text{E}-05$	$7.56935\text{E}-07$	$7.80292\text{E}-08$	$7.53982\text{E}-10$
$2.0\text{E}+19$	$2.16097\text{E}-06$	$6.79243\text{E}-07$	$6.80557\text{E}-08$	$2.87603\text{E}-10$
$5.0\text{E}+19$	$1.59844\text{E}-06$	$5.01530\text{E}-07$	$4.95741\text{E}-08$	$8.04971\text{E}-10$
$1.0\text{E}+20$	$9.69103\text{E}-07$	$3.03930\text{E}-07$	$2.99520\text{E}-08$	$3.77365\text{E}-11$
$2.0\text{E}+20$	$3.56417\text{E}-07$	$1.11747\text{E}-07$	$1.09914\text{E}-08$	$1.15635\text{E}-11$
$5.0\text{E}+20$	$1.77430\text{E}-08$	$5.56222\text{E}-09$	$5.46641\text{E}-10$	$5.21449\text{E}-13$
$1.0\text{E}+21$	$1.19551\text{E}-10$	$3.74776\text{E}-11$	$3.68306\text{E}-12$	$3.49569\text{E}-15$
$2.0\text{E}+21$	$5.42760\text{E}-15$	$1.70148\text{E}-15$	$1.67210\text{E}-16$	$1.58699\text{E}-19$
$5.0\text{E}+21$	$5.07895\text{E}-28$	$1.59218\text{E}-28$	$1.56469\text{E}-29$	$1.48504\text{E}-32$
$1.0\text{E}+22$	$9.79602\text{E}-50$	$3.07092\text{E}-50$	$3.01790\text{E}-51$	$2.86428\text{E}-54$
$2.0\text{E}+22$	$3.64419\text{E}-93$	$1.14240\text{E}-93$	$1.12268\text{E}-94$	$1.06553\text{E}-97$
$5.0\text{E}+22$	$1.87610\text{-}223$	$5.86113\text{-}224$	$5.77979\text{-}225$	$5.48558\text{-}226$

Table B16. Total Photodissociation Rate Coefficients for the Process $\text{CH}_4 + h\nu \rightarrow \text{CH}_3 + \text{H}$ and $\text{CH}_4 + h\nu \rightarrow \text{CH}_2 + \text{H}_2$ for Wavelengths From 116 to 164 nm (Contd)

	Q3 COLUMN DENSITY		Q2 COLUMN DENSITY	
2.0E+13	5.0E+13	1.0E+14	2.0E+14	5.0E+14
1.0E+13	9.08194E-06	9.07579E-06	9.04511E-06	8.98408E-06
2.0E+13	9.08192E-06	9.07577E-06	9.04510E-06	8.98406E-06
5.0E+13	9.08187E-06	9.07573E-06	9.04505E-06	8.98401E-06
1.0E+14	9.08179E-06	9.07564E-06	9.04497E-06	8.98393E-06
2.0E+14	9.08163E-06	9.07548E-06	9.04480E-06	8.98377E-06
5.0E+14	9.08113E-06	9.07499E-06	9.04431E-06	8.98328E-06
1.0E+15	9.08031E-06	9.07417E-06	9.04349E-06	8.98246E-06
2.0E+15	9.07869E-06	9.07254E-06	9.04187E-06	8.98085E-06
5.0E+15	9.07390E-06	9.06775E-06	9.03709E-06	8.97609E-06
1.0E+16	9.06620E-06	9.02941E-06	8.96844E-06	8.86774E-06
2.0E+16	9.05178E-06	9.04565E-06	9.03543E-06	8.95411E-06
5.0E+16	9.01504E-06	9.0892E-06	8.99873E-06	8.9762E-06
1.0E+17	8.96886E-06	8.96276E-06	8.93232E-06	8.87176E-06
2.0E+17	8.90571E-06	8.89964E-06	8.88953E-06	8.80909E-06
5.0E+17	8.78855E-06	8.78254E-06	8.75255E-06	8.69287E-06
1.0E+18	8.65430E-06	8.63847E-06	8.61872E-06	8.46238E-06
2.0E+18	8.45984E-06	8.45401E-06	8.42492E-06	8.36703E-06
5.0E+18	8.07591E-06	8.07037E-06	8.06105E-06	8.04245E-06
1.0E+19	7.62587E-06	7.62057E-06	7.61175E-06	7.59414E-06
2.0E+19	6.88190E-06	6.87711E-06	6.86914E-06	6.85324E-06
5.0E+19	5.09443E-06	5.09088E-06	5.08498E-06	5.07320E-06
1.0E+20	3.08932E-06	3.08717E-06	3.08359E-06	3.07644E-06
2.0E+20	1.13635E-06	1.13556E-06	1.13424E-06	1.13162E-06
5.0E+20	5.65726E-08	5.65333E-08	5.64677E-08	5.63359E-08
1.0E+21	3.81183E-10	3.80917E-10	3.80476E-10	3.79594E-10
2.0E+21	1.73057E-14	1.72936E-14	1.72735E-14	1.71140E-14
5.0E+21	1.61940E-27	1.61827E-27	1.61265E-27	1.60146E-27
1.0E+22	3.12341E-49	3.12124E-49	3.11762E-49	3.11040E-49
2.0E+22	1.16193E-92	1.16113E-92	1.15978E-92	1.15709E-92
5.0E+22	5.98187-223	5.97771-223	5.97078-223	5.95594-223

Table B16. Total Photodissociation Rate Coefficients for the Process $\text{CH}_4 + h\nu \rightarrow \text{CH}_3 + \text{H}$ and $\text{CH}_4 + h\nu \rightarrow \text{CH}_2 + \text{H}_2$ for Wavelengths From 116 to 164 nm

03 COLUMN DENSITY									
1.0E+10	2.0E+10	5.0E+10	1.0E+11	2.0E+11	5.0E+11	1.0E+12	2.0E+12	5.0E+12	1.0E+13
9.08604E-06	9.08604E-06	9.08603E-06	9.08602E-06	9.08600E-06	9.08594E-06	9.08583E-06	9.08563E-06	9.08530E-06	9.08399E-06
2.0E+13	9.08602E-06	9.08602E-06	9.08601E-06	9.08600E-06	9.08598E-06	9.08582E-06	9.08561E-06	9.08500E-06	9.08397E-06
5.0E+13	9.08597E-06	9.08597E-06	9.08596E-06	9.08595E-06	9.08593E-06	9.08587E-06	9.08577E-06	9.08556E-06	9.08495E-06
1.0E+14	9.08589E-06	9.08589E-06	9.08588E-06	9.08587E-06	9.08585E-06	9.08579E-06	9.08569E-06	9.08548E-06	9.08384E-06
2.0E+14	9.08572E-06	9.08572E-06	9.08571E-06	9.08572E-06	9.08569E-06	9.08552E-06	9.08532E-06	9.08470E-06	9.08368E-06
5.0E+14	9.08523E-06	9.08523E-06	9.08522E-06	9.08521E-06	9.08519E-06	9.08513E-06	9.08503E-06	9.08482E-06	9.08421E-06
1.0E+15	9.08441E-06	9.08441E-06	9.08440E-06	9.08439E-06	9.08437E-06	9.08431E-06	9.08421E-06	9.08400E-06	9.08339E-06
2.0E+15	9.08279E-06	9.08278E-06	9.08277E-06	9.08275E-06	9.08269E-06	9.08258E-06	9.08238E-06	9.08176E-06	9.08074E-06
5.0E+15	9.07799E-06	9.07799E-06	9.07798E-06	9.07796E-06	9.07795E-06	9.0779E-06	9.0779E-06	9.07697E-06	9.07595E-06
1.0E+16	9.07029E-06	9.07029E-06	9.07028E-06	9.07027E-06	9.07025E-06	9.07019E-06	9.07009E-06	9.06988E-06	9.06927E-06
2.0E+16	9.05587E-06	9.05587E-06	9.05587E-06	9.05586E-06	9.05584E-06	9.05577E-06	9.05567E-06	9.05547E-06	9.05455E-06
5.0E+16	9.01912E-06	9.01912E-06	9.01912E-06	9.01911E-06	9.01909E-06	9.01902E-06	9.01892E-06	9.01872E-06	9.01708E-06
1.0E+17	8.97293E-06	8.97292E-06	8.97292E-06	8.97291E-06	8.97289E-06	8.97283E-06	8.97273E-06	8.97252E-06	8.97191E-06
2.0E+17	8.90976E-06	8.90975E-06	8.90975E-06	8.90974E-06	8.90972E-06	8.90966E-06	8.90956E-06	8.90935E-06	8.90875E-06
5.0E+17	8.79256E-06	8.79255E-06	8.79255E-06	8.79254E-06	8.79252E-06	8.79246E-06	8.79236E-06	8.79216E-06	8.79156E-06
1.0E+18	8.65826E-06	8.65826E-06	8.65825E-06	8.65824E-06	8.65822E-06	8.65816E-06	8.65806E-06	8.65787E-06	8.65628E-06
2.0E+18	8.46373E-06	8.46373E-06	8.46372E-06	8.46371E-06	8.46369E-06	8.46363E-06	8.46353E-06	8.46334E-06	8.46276E-06
5.0E+18	8.07970E-06	8.07970E-06	8.07969E-06	8.07968E-06	8.07966E-06	8.07961E-06	8.07952E-06	8.07877E-06	8.07784E-06
1.0E+19	7.62294E-06	7.62294E-06	7.62293E-06	7.62293E-06	7.62293E-06	7.62293E-06	7.62293E-06	7.62292E-06	7.62276E-06
2.0E+19	6.88509E-06	6.88508E-06	6.88508E-06	6.88507E-06	6.88505E-06	6.88503E-06	6.88493E-06	6.88429E-06	6.88349E-06
5.0E+19	5.09679E-06	5.09679E-06	5.09679E-06	5.09678E-06	5.09678E-06	5.09677E-06	5.09667E-06	5.09655E-06	5.09561E-06
1.0E+20	3.09075E-06	3.09075E-06	3.09075E-06	3.09074E-06	3.09074E-06	3.09072E-06	3.09068E-06	3.09061E-06	3.09003E-06
2.0E+20	1.13688E-06	1.13661E-06							
5.0E+20	5.65989E-08	5.65989E-08	5.65988E-08	5.65988E-08	5.65988E-08	5.65988E-08	5.65988E-08	5.659823E-08	5.65858E-08
1.0E+21	3.81359E-10	3.81359E-10	3.81359E-10	3.81359E-10	3.81358E-10	3.81358E-10	3.81351E-10	3.81342E-10	3.81271E-10
2.0E+21	1.73137E-14	1.73137E-14	1.73137E-14	1.73137E-14	1.73136E-14	1.73135E-14	1.73133E-14	1.73117E-14	1.73097E-14
5.0E+21	1.62015E-27	1.62015E-27	1.62015E-27	1.62015E-27	1.62013E-27	1.62011E-27	1.62007E-27	1.61996E-27	1.61977E-27
1.0E+22	3.12486E-49	3.12486E-49	3.12486E-49	3.12486E-49	3.12483E-49	3.12479E-49	3.12450E-49	3.12414E-49	3.12414E-49
2.0E+22	1.16247E-92	1.16247E-92	1.16247E-92	1.16247E-92	1.16246E-92	1.16246E-92	1.16245E-92	1.16220E-92	1.16220E-92
5.0E+22	5.98464-223	5.98464-223	5.98464-223	5.98464-223	5.98463-223	5.98462-223	5.98457-223	5.98451-223	5.98395-223

Table B15. Total Photodissociation Rate Coefficients for the Process $\text{HNO}_3 + h\nu \rightarrow \text{OH} + \text{NO}_2$ for Wavelengths From 192 to 325 nm (Contd)

O3 COLUMN DENSITY			
5.0E+18	1.0E+19	2.0E+19	
1.0E+17	1.13617E-05	2.63273E-06	5.63548E-07
2.0E+17	1.13616E-05	2.63273E-06	5.63547E-07
5.0E+17	1.13616E-05	2.63272E-06	5.63547E-07
1.0E+18	1.13615E-05	2.63270E-06	5.63546E-07
2.0E+18	1.13613E-05	2.63266E-06	5.63545E-07
5.0E+18	1.13606E-05	2.63255E-06	5.63541E-07
1.0E+19	1.13596E-05	2.63236E-06	5.63535E-07
2.0E+19	1.13575E-05	2.63199E-06	5.63523E-07
5.0E+19	1.13512E-05	2.63089E-06	5.63486E-07
1.0E+20	1.13409E-05	2.62909E-06	5.63426E-07
2.0E+20	1.13210E-05	2.62559E-06	5.63309E-07
5.0E+20	1.13110E-05	2.62442E-06	5.63288E-07
1.0E+21	1.12486E-05	2.61386E-06	5.62947E-07
2.0E+21	1.11479E-05	2.59721E-06	5.62421E-07
5.0E+21	1.09320E-05	2.56332E-06	5.61404E-07
1.0E+22	1.04726E-05	2.48729E-06	5.58976E-07

O2 COLUMN DENSITY

Table B20. Total Photodissociation Rate Coefficients for the Process $\text{HO}_2 + h\nu \rightarrow \text{O}_2 + \text{H}$ for Wave-lengths From 180 to 274 nm (Contd)

	03 COLUMN DENSITY							
	5.0E+18	1.0E+19	2.0E+19	5.0E+19	1.0E+20	2.0E+20	5.0E+20	1.0E+21
1.0E+17	2.07907E-05	2.61694E-06	6.76290E-08	2.73147E-12	1.93021E-19	1.13611E-33	2.83493E-76	3.46592E-147
2.0E+17	2.07901E-05	2.61691E-06	6.76288E-08	2.73146E-12	1.93020E-19	1.13610E-33	2.83492E-76	3.46592E-147
5.0E+17	2.07884E-05	2.61683E-06	6.76281E-08	2.73145E-12	1.93019E-19	1.13610E-33	2.83491E-76	3.46590E-147
1.0E+18	2.07857E-05	2.61671E-06	6.76269E-08	2.73142E-12	1.93017E-19	1.13609E-33	2.83489E-76	3.46587E-147
2.0E+18	2.07810E-05	2.61648E-06	6.76246E-08	2.73136E-12	1.93014E-19	1.13607E-33	2.83484E-76	3.46582E-147
5.0E+18	2.07695E-05	2.61584E-06	6.76178E-08	2.73119E-12	1.93003E-19	1.13601E-33	2.83471E-76	3.46565E-147
1.0E+19	2.07543E-05	2.61491E-06	6.76067E-08	2.73091E-12	1.92986E-19	1.13592E-33	2.83448E-76	3.46537E-147
2.0E+19	2.07301E-05	2.61326E-06	6.75849E-08	2.73035E-12	1.92951E-19	1.13573E-33	2.83403E-76	3.46482E-147
5.0E+19	2.06766E-05	2.60908E-06	6.75224E-08	2.72868E-12	1.92847E-19	1.13517E-33	2.83268E-76	3.46316E-147
1.0E+20	2.06103E-05	2.60322E-06	6.74242E-08	2.72592E-12	1.92674E-19	1.13425E-33	2.83043E-76	3.46040E-147
2.0E+20	2.05089E-05	2.59329E-06	6.72410E-08	2.72054E-12	1.92334E-19	1.13240E-33	2.82596E-76	3.45491E-147
5.0E+20	2.04746E-05	2.59002E-06	6.71908E-08	2.71987E-12	1.92320E-19	1.13240E-33	2.82596E-76	3.45491E-147
1.0E+21	2.02330E-05	2.56402E-06	6.66700E-08	2.70405E-12	1.91321E-19	1.12696E-33	2.81276E-76	3.43872E-147
2.0E+21	1.99118E-05	2.52725E-06	6.58848E-08	2.67907E-12	1.89725E-19	1.11818E-33	2.79132E-76	3.41247E-147
5.0E+21	1.92876E-05	2.45557E-06	6.43467E-08	2.63067E-12	1.86663E-19	1.10128E-33	2.74998E-76	3.36199E-147
1.0E+22	1.81656E-05	2.31546E-06	6.09623E-08	2.50741E-12	1.78360E-19	1.05394E-33	2.63347E-76	3.21992E-147
2.0E+22	1.64708E-05	2.10592E-06	5.58782E-08	2.32054E-12	1.65729E-19	9.811483E-34	2.45436E-76	3.00131E-147
5.0E+22	1.29553E-05	1.69200E-06	4.63024E-08	1.98438E-12	1.43372E-19	8.53578E-34	2.13714E-76	2.61328E-147
1.0E+23	8.11202E-06	1.06391E-06	2.96259E-08	1.29948E-12	9.486660E-20	5.67955E-34	1.423349E-76	1.73821E-147
2.0E+23	3.52091E-06	4.80146E-07	1.41310E-08	6.52939E-13	4.85590E-20	2.93040E-34	7.34125E-77	8.32785E-48
5.0E+23	5.60412E-07	9.47195E-08	3.31736E-09	1.71564E-13	1.31614E-20	8.03035E-35	2.00348E-77	2.41116E-148
1.0E+24	1.09755E-08	1.96109E-09	7.17647E-11	3.82061E-15	2.95631E-22	1.80327E-36	4.40430E-79	5.10966E-150
2.0E+24	8.94604E-10	1.68185E-10	6.7187E-12	3.23815E-16	2.47282E-23	1.48928E-37	3.56769E-80	4.02630E-51
5.0E+24	8.94366E-10	1.68179E-10	6.17183E-12	3.23815E-16	2.47282E-23	1.48928E-37	3.56769E-80	4.02630E-51
1.0E+25	8.94366E-10	1.68179E-10	6.17183E-12	3.23815E-16	2.47282E-23	1.48928E-37	3.56769E-80	4.02630E-51
2.0E+25	8.94366E-10	1.68179E-10	6.17183E-12	3.23815E-16	2.47282E-23	1.48928E-37	3.56769E-80	4.02630E-51

Table B21. EUV Solar Fluxes and Cross Sections

WAVELENGTH A	PHOTON FLUX $\times 10^{-4}$ PH/CM ² /SEC	$\lambda 10^{-4}$	$\lambda 10^{-4}$		$\lambda 10^{-4}$		$\lambda 10^{-4}$		$\lambda 10^{-4}$	
			N2	O2	N2	O2	N2	O2	N2	O2
1215.7	251.4000	.0100	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000
1025.7	2.6670	1.5800	.0005	0.0000	.9800	0.0000	0.0000	0.0000	0.0000	0.0000
991.0	.5994	1.6000	1.5000	0.0000	1.1000	0.0000	0.0000	0.0000	0.0000	0.0000
977.0	4.4070	4.0000	.7000	0.0000	2.5000	0.0000	0.0000	0.0000	0.0000	0.0000
972.5	.6066	32.0000	36.0000	0.0000	25.0000	0.0000	0.0000	0.0000	0.0000	0.0000
949.7	.3044	6.3000	5.2000	0.0000	4.0000	0.0000	0.0000	0.0000	0.0000	0.0000
944.5	.1340	3.2000	1.0000	0.0000	2.6000	0.0000	0.0000	0.0000	0.0000	0.0000
937.8	.1635	5.0000	10.0000	0.0000	3.0000	0.0000	0.0000	0.0000	0.0000	0.0000
933.4	.1433	18.0000	.2000	0.0000	8.0000	0.0000	0.0000	0.0000	0.0000	0.0000
930.7	.2631	26.0000	4.8000	0.0000	17.0000	0.0000	0.0000	0.0000	0.0000	0.0000
926.2	.1129	6.7000	2.0000	0.0000	6.7000	0.0000	0.0000	0.0000	0.0000	0.0000
	.0284	7.4000	100.0000	0.0000	3.3000	0.0000	0.0000	0.0000	0.0000	0.0000
	.1134	7.4000	50.0000	0.0000	3.3000	0.0000	0.0000	0.0000	0.0000	0.0000
	.3970	7.4000	10.0000	0.0000	3.3000	0.0000	0.0000	0.0000	0.0000	0.0000
	.5388	7.4000	2.0000	0.0000	3.3000	0.0000	0.0000	0.0000	0.0000	0.0000
890-911	.3403	7.4000	.2000	0.0000	3.3000	0.0000	0.0000	0.0000	0.0000	0.0000
	1.1064	9.3000	10.0000	3.0000	4.5000	0.0000	3.0000	0.0000	6.1000	0.0000
904	2.5816	9.3000	2.0000	3.0000	4.5000	0.0000	3.0000	0.0000	6.1000	0.0000
	.3625	11.0000	6.3000	3.0000	6.3000	0.0000	3.0000	0.0000	6.1200	0.0000
860-890	1.1852	7.0000	10.0000	2.9000	4.0000	0.0000	2.9000	5.8000	0.0000	0.0000
	1.7778	7.0000	2.0000	2.9000	4.0000	0.0000	2.9000	5.8000	0.0000	0.0000
830- 60	.5436	9.0000	10.0000	2.8000	4.0000	0.0000	2.8000	5.2000	0.0000	0.0000
	.8154	9.0000	2.0000	2.8000	4.0000	0.0000	2.8000	5.2000	0.0000	0.0000
	.0950	10.0000	20.0000	2.6000	3.7000	0.0000	2.6000	5.0000	0.0000	0.0000
	.0950	10.0000	41.0000	2.6000	4.0000	0.0000	2.6000	5.8000	0.0000	0.0000
	.0950	10.0000	7.5000	2.6000	4.0000	0.0000	2.6000	5.0000	0.0000	0.0000
	.0950	13.0000	6.0000	2.6000	5.1000	0.0000	2.6000	5.0000	0.0000	0.0000
	.0950	15.0000	2.3000	2.6000	6.0000	0.0000	2.6000	5.0000	0.0000	0.0000
	.0950	26.0000	7.0000	2.6000	10.0000	0.0000	2.6000	5.0000	0.0000	0.0000
	.0950	26.0000	2.1000	2.6000	10.0000	0.0000	2.6000	5.0000	0.0000	0.0000
	.0654	19.0000	50.0000	2.9000	7.0000	0.0000	2.900	4.8000	0.0000	0.0000
	.1962	19.0000	10.0000	2.9000	7.0000	0.0000	2.900	4.8000	0.0000	0.0000
	.3924	19.0000	2.0000	2.9000	7.0000	0.0000	2.900	4.8000	0.0000	0.0000
	.0027	26.0000	100.0000	3.0000	10.0000	65.0000	3.0000	4.3000	0.0000	0.0000
	.0539	26.0000	50.0000	3.0000	10.0000	33.0000	3.0000	4.3000	0.0000	0.0000
	.2129	26.0000	25.0000	3.0000	10.0000	16.0000	3.0000	4.3000	0.0000	0.0000
	.3814	28.0000	25.0000	2.9000	10.0000	11.0000	2.900	4.4000	0.0000	0.0000
	.2107	24.0000	10.0000	2.9000	13.0000	8.0000	2.900	4.3000	0.0000	0.0000
	.1103	23.0000	44.0000	2.9000	7.0000	33.0000	2.900	4.3000	0.0000	0.0000
	.786.5									
	.780.3	.2289	28.0000	30.0000	2.0000	11.0000	17.0000	2.900	4.3000	0.0000
	.790.1	.2869	18.0000	12.0000	7.0000	11.0000	7.8000	7.000	4.1000	0.0000
	.787.7	.2107	24.0000	10.0000	2.9000	11.0000	35.0000	3.300	3.9000	0.0000
	.740-770	.0128	19.0000	50.0000	3.3000	11.0000	17.0000	3.300	3.9000	0.0000
	.1698	19.0000	24.0000	3.3000	11.0000	51.0000	3.000	4.0000	0.0000	0.0000
	.765.1	.1659	23.0000	67.0000	3.0000	12.0000	22.0000	2.900	3.9000	0.0000
	.760.4	.0934	20.0000	40.0000	2.9000	10.0000	22.0000	2.900	3.9000	0.0000
	.710-740	.0895	30.0000	23.0000	5.5000	21.0000	5.5000	5.500	3.5000	0.0000

Table B21. EUV Solar Fluxes and Cross Sections (Contd)

WAVELENGTH Å	PHOTON FLUX $\times 10^9$ PH/CM ⁻² /SEC	O^2 $\times 10^{-4}$	N_2 $\times 10^{-4}$	HE $\times 10^{-10}$	
				O $\times 10^{-4}$	N_2 $\times 10^{-4}$
680-710	.1290	24.0000	23.0000	6.5000	20.0000
703	.0673	26.0000	23.0000	6.5000	21.0000
630-680	.0004	22.0000	50.0000	6.0000	50.0000
	.0627	22.0000	24.0000	6.0000	19.0000
629.7	1.5203	30.0000	23.0000	9.0000	24.0000
625.3	.2103	25.0000	24.0000	9.0000	24.0000
609.8	.5302	27.0000	24.0000	9.0000	25.0000
599.6	.1600	28.0000	23.0000	9.0000	27.0000
584.3	1.2700	23.0000	23.0000	9.5000	23.0000
554	.7200	26.0000	25.0000	9.5000	24.0000
521.0	.0600	21.0000	24.0000	9.3000	23.0000
507	.1600	23.0000	23.0000	9.3000	22.0000
425-504	.4429	22.0000	23.0000	9.5000	22.0000
499.3	.1139	22.0000	23.0000	9.5000	22.0000
465.2	.2724	21.0000	23.0000	10.0000	21.0000
417	.0782	20.0000	21.0000	10.5000	20.0000
370-460	.1715	20.0000	20.0000	10.0000	20.0000
368.1	.6500	18.0000	17.0000	9.5000	18.0000
365	.1500	18.0000	16.0000	9.5000	18.0000
360.7	.3200	17.0000	16.0000	9.5000	17.0000
335.4	.1600	17.0000	14.0000	9.0000	17.0000
303.8	7.7000	17.0000	12.0000	8.5000	17.0000
284.1	.2100	15.0000	11.0000	8.5000	15.0000
270-370	1.1770	16.0000	13.0000	8.5000	18.0000
230-270	2.3960	13.0000	10.0000	7.5000	22.0000
205-230	.6035	10.0000	7.0000	6.5000	20.0000
176-205	2.0881	6.4000	5.6000	4.0000	12.8000
153-176	1.2133	5.2000	4.4000	3.0000	10.4000
100-153	.2017	3.0000	2.7000	1.7000	7.5000
90-100	.0942	2.0000	1.2000	1.0000	6.0000
80-90	.1070	1.5000	.8900	.7500	4.5000
76.0	.0167	1.1000	.6700	.5500	4.4000
70-80	.0648	1.1000	.6500	.5500	4.4000
60-70	.0742	.8000	.4500	.3900	4.0000
50-60	.0643	.5100	.3000	.2500	3.1000
40-50	.0284	.3100	.1800	.1600	2.2000
10-40	.0040	.1500	.0900	.0800	1.5000
9	0.0000	.1000	.0650	.0500	3.9000
7	0.0000	.0520	.0300	.0260	2.6000
5	0.0000	.0200	.0120	.0100	1.4000
3	0.0000	.0046	.0028	.0023	.5400
1	0.0000	.0002	.0001	.0001	.0350

Table B22. Electron Impact Ionization Cross Sections

PROCESS	I	$c_0 f_0$	Ω	ν	γ	P
$e + N_2 \rightarrow N + N^+ + 2e$	25.0	0.380	0.96	2.0	1.0	1.2
$e + O_2 \rightarrow O + O^+({}^4S) + 2e$	18.0	0.400	0.93	3.0	1.0	1.1
$e + O_2 \rightarrow O + O^+({}^2D) + 2e$	22.0	0.250	0.93	3.0	1.0	1.1
$e + O \rightarrow O^+({}^4S) + 2e$	14.0	0.290	0.85	1.0	0.3	1.2
$e + O \rightarrow O^+({}^2D) + 2e$	17.0	0.360	0.85	1.0	0.3	1.2
$e + N_2 \rightarrow N_2^+ + 2e$	15.0	0.370	0.80	3.0	1.0	1.2
$e + O_2 \rightarrow O_2^+ + 2e$	12.0	0.058	0.80	2.0	1.0	1.1

Table B23. Rate Coefficients for Energetic Electron Impact Reactions

1. $N_2 + e \rightarrow N + N^+ + 2e$				
z(km)	Solar Zenith Angle			
	49.78°	61.50°	76.22°	97.16°
95	0	0	0	0
100	3.54E-12	7.83E-13	3.37E-15	0
130	7.46E-10	5.79E-10	3.07E-10	0
170	2.60E-09	2.25E-09	1.45E-09	0
210	3.92E-09	3.69E-09	3.03E-09	0
250	4.73E-09	4.61E-09	4.22E-09	0
300	5.27E-09	5.21E-09	5.04E-09	0
400	2.63E-09	2.60E-09	2.52E-09	0

2. $O_2 + e \rightarrow O + O^+({}^4S) + 2e$				
z(km)	Solar Zenith Angle			
	49.78°	61.50°	76.22°	97.16°
95	0	0	0	0
100	4.38E-12	9.68E-13	4.16E-15	0
130	1.00E-09	7.49E-10	3.87E-10	0
170	4.60E-09	3.84E-09	2.20E-09	0
210	7.31E-09	6.81E-09	5.35E-09	0
250	8.86E-09	8.59E-09	7.73E-09	0
300	9.76E-09	9.61E-09	9.25E-09	0
400	4.38E-09	4.81E-09	4.62E-09	0

Table B23. Rate Coefficients for Energetic Electron Impact Reactions (Contd)

3.	$O_2 + e \rightarrow O + O^+(^2D) + 2e$				
	$z(km)$	Solar Zenith Angle			
		49.78°	61.50°	76.22°	97.16°
95	0	0	0	0	0
100	2.06E-12	4.56E-13	1.97E-15	0	0
130	4.11E-10	3.22E-10	1.74E-10	0	0
170	1.35E-09	1.17E-09	7.74E-10	0	0
210	2.01E-09	1.90E-09	1.57E-09	0	0
250	2.42E-09	2.36E-09	2.17E-09	0	0
300	2.70E-09	2.67E-09	2.59E-09	0	0
400	1.35E-09	1.33E-09	1.29E-09	0	0
4.	$O - e \rightarrow O^+(^4S) + 2e$				
	$z(km)$	Solar Zenith Angle			
		49.78°	61.50°	76.22°	97.16°
95	0	0	0	0	0
100	5.90E-12	1.29E-12	5.53E-15	0	0
130	2.08E-09	1.29E-09	5.73E-10	0	0
170	2.37E-09	1.79E-09	7.26E-09	0	0
210	4.26E-08	3.85E-08	2.70E-08	0	0
250	5.09E-08	4.88E-08	4.22E-08	0	0
300	5.22E-08	5.08E-08	4.82E-08	0	0
400	2.61E-08	2.54E-08	2.41E-08	0	0
5.	$O + e \rightarrow O^+(^2D) + 2e$				
	$z(km)$	Solar Zenith Angle			
		49.78°	61.50°	76.22°	97.16°
95	0	0	0	0	0
100	4.23E-12	9.30E-13	3.98E-15	0	0
130	1.28E-09	8.51E-10	3.40E-10	0	0
170	1.07E-08	8.35E-09	3.80E-09	0	0
210	1.84E-08	1.68E-08	1.22E-08	0	0
250	2.22E-08	2.13E-08	1.86E-08	0	0
300	2.36E-08	2.31E-08	2.20E-08	0	0
400	1.18E-08	1.15E-08	1.10E-08	0	0

Table B23. Rate Coefficients for Energetic Electron Impact Reactions (Contd)

6. $N_2 + e \rightarrow N_2^+ + 2e$				
z(km)		Solar Zenith Angle		
	49.78°	61.50°	76.22°	97.16°
95	0	0	0	0
100	6.12E-12	1.34E-12	5.80E-15	0
130	1.55E-09	1.11E-09	5.52E-10	0
170	9.31E-09	7.51E-09	3.86E-09	0
210	1.54E-08	1.42E-08	1.07E-08	0
250	1.87E-08	1.80E-08	1.60E-08	0
300	2.02E-08	1.98E-08	1.90E-08	0
400	1.01E-08	9.91E-09	9.49E-09	0
7. $O_2 + e \rightarrow O_2^+ + 2e$				
z(km)		Solar Zenith Angle		
	49.78°	61.50°	76.22°	97.16°
95	0	0	0	0
100	1.81E-12	3.97E-13	1.70E-15	0
130	6.32E-10	3.94E-10	1.76E-10	0
170	7.14E-09	5.40E-09	2.19E-09	0
210	1.29E-08	1.17E-08	8.18E-09	0
250	1.55E-08	1.48E-08	1.28E-08	0
300	1.60E-08	1.53E-08	1.45E-08	0
400	7.90E-09	7.67E-09	7.27E-09	0
8. $N_2 + e \rightarrow N + N(^2D) + e$				
z(km)		Solar Zenith Angle		
	49.78°	61.50°	76.22°	97.16°
95	0	0	0	0
100	1.64E-11	3.59E-12	1.54E-14	0
130	6.27E-09	3.71E-09	1.60E-09	0
170	9.42E-08	6.93E-08	2.53E-08	0
210	1.82E-07	1.63E-07	1.11E-07	0
250	2.22E-07	2.12E-07	1.81E-07	0
300	2.11E-07	2.03E-07	1.91E-07	0
400	1.05E-07	1.01E-07	9.55E-08	0

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Appendix C

The Computer Code

The computer code used to develop the simulated atmosphere described in this report is presented in detail in this Appendix.

To conserve computer memory, the code was written in OVERLAYS. Once all the initialization is completed at the beginning of a run, there is no need to retain in memory all the code that accomplished this.

The code is written in FORTRAN IV for a CDC 6600 mainframe computer. It also utilizes a large block of extended core storage for saving the concentration arrays, the average velocity arrays, and the reaction rate coefficient arrays at the beginning of each time step. If extended core is not available, these sections of the code have to be rewritten to use whatever disk or, as a last resort, tape storage is available. Using disk or tape storage significantly increases the residence time of the code on the computer.

C1. THE MAIN OVERLAY (0,0)

The main overlay is relatively short. It defines all the blank COMMON variables shared by all the overlays. It also defines the parameters that describe the simulation to be generated as well as the error criteria. It calls OVERLAYS (1,0) and (2,0).

C1.1 Blank COMMON

There are 692 height steps between the lower and the upper boundaries in this model. The 424 height steps between the lower boundary and 120 km represent the region over which the negative ions and the cluster ions are calculated.

N is one of the dependent variable arrays. It contains the altitude distribution of the concentrations of the 26 species in the transport mode plus the arrays for O(¹D), N₂⁺, O(²D), and the electrons.

CONC is also a dependent variable array. It contains the height distributions for those species considered over the smaller altitude region.

U is another dependent variable array. Here are stored the values at each altitude of the average velocities of those species in the transport mode.

NZERO and UZERO are working arrays that contain the values of N and U for the particular species whose set of equations is being solved at any given time.

FORM and REMV contain the altitude arrays of the rate of chemical formation [P in Eq. (1)] and chemical removal [L in Eq. (1)] for the species whose equations are being solved.

E and F contain the altitude arrays of E and F in Eq. (28) for the species under consideration.

Z is the altitude array in kilometers.

DEDY is the array containing the fixed vertical turbulent diffusion coefficients.

TEMP is the array containing the fixed temperature profile.

DTEMDZ is the array of the values

$$\frac{1}{T(z)} \frac{dT(z)}{dz} \quad (C1)$$

DELTAZ is the array of height increments in centimeters.

SUNSET is the array containing the sunset solar zenith angle at each altitude.

GRAV is the sea level acceleration of gravity.

COSD = cos (declination)*cos (latitude).

SIND = sin (declination)*sin (latitude). Both COSD and SIND are used in computing the solar zenith angle.

TIME is the independent variable. Time is measured in seconds, with time = 0 as noontime of solution day 1.

CXI is the solar zenith angle in radians and is a function of the local time.

RADIUS is the radius of the earth in kilometers.

ISPECI is the total number of species included in the code.

KSPECI is the number of species in the transport mode.

NREAC is the number of temperature independent chemical reactions.

NREAC2 is the number of temperature dependent chemical reactions.

K is the number of height steps over which the equations are to be integrated.

KMIN1 = K - 1.

ITURB is the index of the height of the turbopause in the Z array.

CNINOON is the noontime solar zenith angle.

SCXI is the local solar zenith angle in degrees.

C1.2 Labeled COMMON

C1.2.1. COMMON/ESM/

OLDN, OLDCON, and OLDU are the extended core arrays corresponding to N, CONC, and U respectively.

RATES is the extended core array containing the values of the height dependent reaction rate constants for the 88 temperature dependent reactions. These four variables are also defined in the LEVEL statement in all programs where they apply.

C1.2.2 COMMON/TIMES/

DAY\$ is the variable that contains the number of solution days over which the equations will be solved.

DELT always contains the current value of the time step in seconds.

TMAX is the maximum value in seconds that the time step will be allowed to have. This value is 1800 seconds in code presented here.

TSFT is the time in seconds after noontime of sunset at the lower boundary.

TRISE is the time in seconds after noontime of sunrise at the upper boundary.

C1.2.3 COMMON/ALTITUD/

ZBOT is the altitude of the lower boundary in kilometers.

ZTOP is the altitude of the upper boundary in kilometers.

TURB is the altitude of the turbopause in kilometers.

ZTRANS is the altitude of the day-night transition height. During the day, it is set at 1km, and, during the night, it is set at 400 km.

C1.2.4 COMMON/RESTART/

IFRAME is the number given to a particular solution file or the interrupt file Tape 4. These files are numbered sequentially from 1.

TOFF is the number of seconds that must be allowed the program to dump onto the interrupt file all the information needed to restart the code at the point when it was interrupted. The final solution of these sets of equations takes many hours of CDC 3200 computer time. Because of this, the code is normally run for a few hours (usually 4), then stopped, and a dump is taken. Later, the code is restarted at this point, continued for another period of time, and another dump is taken.

This routine is repeated until the calculations reach a point where they repeat the solution over each diurnal cycle (usually about 20 days of problem time).

TYME is the time in seconds of the current computer run. The computer stops executing the code and dumps when

TYME - (time running) < TOFF.

TODAY is the date when the run was made and is printed out as XX/XX/XX.

C1.2.5 COMMON/OUTPUTS/

TIMOUT is a variable that gives the code the frequency in hours with which to send the solutions to the OUTPUT and the interrupt files. The solutions after every time step are not saved. Only those solutions occurring at TIMOUT hours and at interrupts are sent to OUTPUT and are preserved on Tape 4.

OPRINT is the variable containing the frequency in hours at which solutions are to be written on Tape 4. Except on interrupts, TIMOUT is incremented by OPRINT every time a solution set is written to Tape 4.

PRINT is a logical parameter under program control that determines whether writing to Tape 4 will take place after a solution is accepted (PRINT = .TRUE.) or will not (PRINT = .FALSE.).

C1.2.6 COMMON/ERRORS/

ERRMAX is the criterion for acceptance of a solution. When the relative change of all species at all altitudes over successive iterations is less than ERRMAX, the iteration is stopped and the solution for that time step accepted.

ERSPECI is the criterion for stopping the iteration of individual species. Since the solutions converge faster for some species than for others, a good deal of computer time can be saved by no longer iterating the equations for those species whose relative change is less than ERSPFCI. Of course, ERSPFCI must be much smaller than ERRMAX.

ITMAX is the maximum number of iterations to be allowed for any given solution. If any species have not converged after ITMAX iterations, the iterating is stopped and the last solutions computed are accepted. With ITMAX = 100, this situation has never arisen.

C1.2.7 COMMON/CONSTAN/

PI = 3.14159265359.

RADDEG = PI/180 is the conversion factor for degrees to radians.

DEGRAD = 180/PI is the conversion factor for radians to degrees.

RADSEC = PI/43200 is the conversion factor for changing local time in seconds to local hour angle in radians.

BOLTZ = 1.38047×10^{-16} ergs/K is Boltzmann's constant.

SIN2I is the square of the sine of the magnetic dip angle.

ATCON = 1.660356×10^{-24} is the mass of the hydrogen atom in grams.

C1.2.8 COMMON/JAYS/

The labeled COMMON/JAYS/ contains the photodissociation rate coefficients for reactions 192 through 205. The locations reserved in this labeled COMMON are initialized from file Tape 3 by PROGRAM INITIAL.

C1.2.9 COMMON/COLFRE/

The array D contains the invariant part of the collision frequencies ν_{ij} .

$$D[J,I] = \frac{1}{2} (\sigma_I + \sigma_J)^2 \left[2\pi k \frac{(m_I + m_J)}{m_I m_J} \right]^{1/2}. \quad (C2)$$

C1.2.10 COMMON/ATCONS/

The arrays in the labeled COMMON/ATCONS/ all pertain to the atomic and molecular constants.

ATWT contains the atomic or molecular weights of the 26 species in the transport mode.

MASS=ATWT*ATCON contains the mass in grams of the 26 species in the transport mode.

THERM contains the thermal diffusion factors for the 26 species in the transport mode. Where values are not known, the thermal diffusion factor is set to unity, which removes any thermal diffusion effect for that species.

SYMBOL contains the alphanumeric identification of each of the 56 species in the model.

POLAR contains the polarizabilities of the species in the transport mode. For species whose polarizabilities are not known, the value is set to zero to remove the polarizability effect.

C1.2.11 COMMON/RATECON/

The arrays in labeled COMMON/RATECON/ all pertain to the rate coefficients for the chemical reactions.

The arrays AK, BK, and CK contain the A's, B's, and C's of Eq. (A1) for the B height dependent reaction rate coefficients.

The array DK contains the rate coefficients for the 215 chemical reactions at the particular altitude under consideration.

The array CYMB contains the alphanumeric equivalence of the 215 chemical reactions themselves.

C1.2.12 COMMON/IONS/

The arrays in labeled COMMON/IONS/ all pertain to the ionization of the atmospheric gases.

WAVE contains the alphanumeric identification of the lines or wavelength bands in the EUV.

FLUX contains the solar flux in photons/cm²/sec/ \AA at the top of the atmosphere at each of the above lines or in each of the wavelength bands.

SIGMA contains the absorption and the ionization cross sections for O₂, N₂, and O, in that order.

R contains the rate coefficients for the eight energetic electron reactions at the four zenith angles and eight heights listed in Table B22.

The main overlay sets the values of the control parameters before calling OVERLAY (1, 0). Upon return from OVERLAY (1, 0), the initialization of the code for the run to be made is completed. The initial values of all of the dependent variables N, CONC, and U are then saved in extended core storage. Finally, OVERLAY (2, 0) is called and retains control until the calculations are terminated.

C2. OVERLAY (1,0) PROGRAM INITIAL

The purpose of OVERLAY (1, 0) PROGRAM INITIAL is to initialize the code for the particular run to be made. In the dimension statement—

S contains the collision cross sections in cm² of each of the species in the transport mode. This information is required in the computation of the collision frequencies ν_{ij} .

R MIX contains the ground level mixing ratios for the transport species.

SIG contains the photoionization cross sections for atomic nitrogen for the lines and bands given in array WAVE. These cross sections scaled by 10¹⁸ are listed in the DATA statement of this overlay.

Following this DATA statement are the FUNCTION definitions. The first is for the sea level value for the acceleration of gravity as a function of the latitude Φ . It is given by^{C1}

$$\text{GRAV}(\Phi) = 978.0356 [1 + 0.0052835 \sin^2(\Phi) - 5.9 \times 10^{-6} \sin^2(2\Phi)]. \quad (\text{C3})$$

The second is for the radius of the earth as a function of the colatitude and the sea level acceleration of gravity. It is given by^{C1}

C1. List, R.J., Ed. (1963) Smithsonian Meteorological Tables, Sixth Revised Edition, Washington, D.C.

altitude whose index is IALT. First, this subroutine computes the total concentration M as

$$Y(57) = \sum_{i=1}^{56} n_i.$$

Then it moves the rate coefficients for the 215 reactions at the altitude under consideration into the K region. Following this, it ensures that the photodetachment and photodissociation of the negative ions are removed if the code is in the nighttime. The remainder of the code computes the P and L for species KIND and stores them in the appropriate FORM and REMV regions, respectively. The bulk of this subroutine as well as SUBROUTINE CHEMPR were not hand-coded for obvious reasons. To avoid the very tedious task of hand-coding these routines every time the chemistry set or the species set is changed, a set of programs was written to perform the task of writing these codes. These programs can be found in Keneshea.³

3.11 SUBROUTINE RATECN

SUBROUTINE RATECN sets the reaction rate coefficients for all the photo-reactions at the altitude whose index is IALT. The first three DATA statements in this subroutine contain, respectively, the solar fluxes ($\times 10^{-9}$ photons), the ionization cross sections ($\times 10^{18} \text{ cm}^2$), and the effective absorption cross sections ($\times 10^{18} \text{ cm}^2$) for $O_2(^1\Delta g)$ as given by Huffman.²² The next two DATA statements contain the altitudes (km) and the solar zenith angles corresponding to the input arrays of the rate coefficients for the energetic electrons. Following this is the DATA statement specifying the rate coefficient for ionization by cosmic rays. The first executable statement of the program is the FUNCTION definition for the correction at altitude ZZ for the sea level acceleration of gravity.

The code first computes all the rate coefficients for the current altitude and stores them in the DK array. On the initial call to this subroutine, the fluxes and the cross sections given in the DATA statements are scaled to their proper orders of magnitude, and the rate coefficients for the energetic electron processes are replaced by their natural logarithms. Next, the noontime values of H x F for O_2 , N_2 , and O are computed and saved for later use in computing the scattered radiation. Here, H is the scale height of the species and F is the Chapman approximation used in computing the absorption. If the code is in the nighttime, then transfers are made to statements near the end of the code since all photo processes are inactive during the night. If the code is in the daytime, then the photo rate coefficients are determined for the immediate daytime conditions. If the code is in the

where, as before, ℓ and $\ell + 1$ represent values at the beginning and at the end of the time step, respectively. The profile of O(¹D) is computed once from Eq. (C21) each time that this subroutine is called, with P_k^ℓ and L_k^ℓ used instead of $P_k^{\ell+1}$ and $L_k^{\ell+1}$. Also computed in this subroutine are the concentrations of : (1) all the positive ions not in the transport mode; (2) all the negative species; (3) all the positive cluster ions; and (4) HNO₂ and NO₃ between 50 and 120 km.

Eq. (C21) requires the value of the chemistry terms P and L at the end of the time step which in turn require the value of n at the end of the time step. Since this is the quantity being determined, Eq. (C21) will have to be iterated. Upon each iteration, the chemistry terms are computed with the latest values of n. When the solution has converged, the value of n at $\ell + 1$ will have been achieved and the chemistry will have been computed with this value.

At the beginning of each iteration, the value of the concentration at the current altitude is saved temporarily in the array H. The new value is stored in the appropriate cell of the F array. To ensure a more rapid convergence of the solution, the value of the concentration after the current iteration is replaced by the average value.

$$n = \frac{H + F}{2}$$

Then the relative difference

$$\frac{|H - n|}{H}$$

is computed for each species. If the relative change for each species over the current iteration is less than 20 percent, the solution is accepted for that height step and the iterations are started at the next higher altitude. If any species has a relative change greater than 20 percent, all the concentrations are iterated again. This procedure is repeated until all species have relative changes at all altitudes of less than 20 percent. The total number of iterations allowed at each altitude is 100. This number has always been more than sufficient to generate a solution.

When the concentrations of all the required species have been computed between 50 and 120 km, this subroutine then computes the electron concentration between 120 km and the upper boundary from charge balance.

C.13 SUBROUTINE CHEM

SUBROUTINE CHEM computes the quantities P and L for species KIND at the

The variable TIMOUT is used to keep account of the next hour in problem time when a solution will be sent to the output file. The code ensures that this time will occur exactly by resetting the time step if the current time exceeds this time. The RESET and PRINT switches are set to .TRUE., and the value of TIMOUT is incremented by OPRINT to update it for the next printout. When this routine is exited, DELT contains the time increment in seconds over which the next solution will be made, TIME contains the total elapsed problem time in seconds, and CHI contains the solar zenith angle in radians at the end of the current time step.

C3.12 SUBROUTINE CHEMION

SUBROUTINE CHEMION computes the concentrations of all species in the chemistry mode. This subroutine also computes the largest negative species from charge balance. Below 120 km, this could be a negative ion or the electrons. Above 120 km, it is always the electrons.

If the divergence of the flux in Eq. (1) is set to zero, then that equation can be written as

$$\frac{dn_i}{dt} = P_i - n_i L_i . \quad (C19)$$

Two convenient choices exist for solving these ordinary differential equations. They can be solved using the Kutta-Merson numerical integration scheme available in SUBROUTINE ODE, or they can be solved using simple finite differences. Since much more computer time is required to solve the equations by the Kutta-Merson scheme and since it is generally a minor species whose concentration is being computed, the more economical finite differences are used to solve Eq. (C19). Writing Eq. (C19) in finite difference notation at level k gives

$$\frac{n_k^{\lambda+1} - n_k}{\Delta t} = P_k^{\lambda+1} - n_k^{\lambda+1} L_k^{\lambda+1} . \quad (C20)$$

The concentration $n_k^{\lambda+1}$ at the end of the time step is then

$$n_k^{\lambda+1} = \frac{n_k^\lambda + \Delta t P_k^{\lambda+1}}{1 + \Delta t L_k^{\lambda+1}} \quad (C21)$$

C3.8 SUBROUTINE ODE

SUBROUTINE ODE solves a set of N ordinary differential equations using the Kutta-Merson numerical integration scheme. This program is explained in detail elsewhere.² The solution is taken over the altitude range DZ starting at altitude SZ. The solutions are returned in array START.

C3.9 SUBROUTINE RHS

SUBROUTINE RHS is called by SUBROUTINE ODE to supply to ODE the values of the right hand side of Eq. (42). The values are computed at altitude ZE using the intermediate values v computed in SUBROUTINE ODE. The results of the evaluation are returned in the array SLOPE. The intermediate values at altitude ZE are computed using the linear interpolation coefficients derived in SUBROUTINE VELOCITY.

C3.10 SUBROUTINE COLFRE

SUBROUTINE COLFRE computes the sum of the collision frequencies (SNU) of species i with all the other species at the altitude whose index is j. It also computes the quantities $\sum_j \nu_{ij} w_j$ required in Eq. (22) for species i. In addition, it sets the values of omega (OMEG) in Eq. (17). OMEG is computed from Eq. (10) if i represents an ion, and is set equal to unity if i represents a neutral species.

C3.11 SUBROUTINE TIMER

Except for doubling of the time step after a successful integration and halving of the time step on an error, all changes to the time and the time step are done in SUBROUTINE TIMER. This subroutine decides whether or not the next integration will end at a print step. It also determines if the next time step will be in the daytime, nighttime, or in morning or evening twilight. The code advances normally in time until this subroutine determines that the sun is about to rise at the upper boundary or is about to set at the lower boundary. When either of these situations is about to happen, this subroutine adjusts the time step so that the next integration will end at the beginning of morning or evening twilight. During these twilight periods, the concentrations of some of the atmospheric species change rapidly. These changes force the code to reduce the time step to successively smaller and smaller values before a solution can be accepted. To avoid the large amount of computer time required to accomplish this, the code automatically reduces the time step to 15 seconds with the onset of morning or evening twilight. The LOCAL switch SET controls this. The time increment is then allowed to double in PROGRAM INTEG after each successful integration until it reaches TMAX.

$$\frac{n_j + 1}{n_j} < 100.$$

If the ratio $n_{j+1}/n_j > 100$, the velocity is set to zero at that altitude.

C3.7 SUBROUTINE VELOCITY

SUBROUTINE VELOCITY sets up the parameters required for computing the upper boundary value of the velocity of species i. In order to use any numerical integration scheme to solve Eq. (42) for the velocity w at the upper boundary, some means of interpolating values of n, P, L, and dn/dz between the upper boundary and one step below must be provided. What is used here is a simple linear interpolation:

$$\begin{cases} n_K = az_K + b \\ n_{K-1} = az_{K-1} + b \end{cases} \quad (C17)$$

Solving these equations for a and b gives

$$a = \frac{n_K - n_{K-1}}{\Delta z_K}.$$

and

$$b = \frac{n_{K-1} z_K - n_K z_{K-1}}{\Delta z_K}. \quad (C18)$$

In the function definitions at the beginning of SUBROUTINE VELOCITY, AF is the definition of a and BF is the definition of b. The subroutine then proceeds to compute the a and b coefficients for the equations n(AN,BN), P(AP,BP), and L(AL,GL). These are stored in labeled COMMON/RHSIDE/ for use by SUBROUTINE RHSINTER. This subroutine then calls SUBROUTINE ODE to solve the ordinary differential equations Eq. (42) for species i over the last height step.

the total concentration

$$N = \sum_{i=1}^{26} n_i,$$

the mean molecular weight in grams of the atmosphere

$$\bar{m} = \frac{\rho}{N}, \quad (C15)$$

and the mean mass velocity

$$w_o = \frac{\sum_{i=1}^{26} \rho_i w_i}{\sum_{i=1}^{26} \rho_i} \quad (C16)$$

The negative ions and the positive cluster ions, which are always very minor species, are excluded from the above sums.

C3.6 SUBROUTINE EFORU

For the species whose number is ICF, SUBROUTINE EFORU computes its E and F arrays of Eq. (28) if ICF is positive, or its average velocity array from Eq. (18) if ICF is negative.

The computation of the E and F arrays is fairly straightforward. At each altitude, the A(AJ), B(BJ), C(CJ), and D(DJ) values of Eq. (21) are computed. Then the E and F arrays are computed with the upper and lower boundary conditions imposed.

The specific velocities are also calculated in this subroutine since their evaluation requires much of the same coding as do the E and F arrays. Both computations call SUBROUTINE VELOCITY to obtain the upper boundary velocity for those species that are not assumed to be in diffusive equilibrium there. To avoid unrealistically large values of the velocities when the concentration is very small, the concentration of each species n_j and the ratio of n_{j+1}/n_j are investigated at each altitude. If $n_j > 1.0 \times 10^{-10} \text{ cm}^{-3}$, then its velocity at altitude j is computed. For values of $n_j < 1.0 \times 10^{-10}$, the velocity is computed if the ratio

Performing the integration results in

$$n = n_K \exp \left\{ - \frac{mg_O}{kT_K} \left[\frac{(z - z_K)}{\left(\frac{z_K}{R} + 1 \right) \left(\frac{z}{R} + 1 \right)} \right] \right\} \quad (C12)$$

where the subscript K refers to quantities at the upper boundary. After the concentrations have been extrapolated to 500 km, the actual smoothing routine is called, where the profile is smoothed with a 7-point running mean. At this point, the code continues to loop back until all the required smoothing is accomplished. If no ion profile was smoothed, the subroutine is exited here. If at least one of the smoothed profiles was a positive ion profile, then the negative species must be readjusted to ensure charge neutrality in the atmosphere. Below 120 km, the largest negative species is recomputed from charge balance.

$$n_i = \sum_j n_j^+ - \sum_{j \neq i} n_j^- \quad (C13)$$

Above 120 km the electron concentration is recomputed from

$$n_e = \sum_j n_j^+$$

C.3.4 SUBROUTINE SMOOTH

SUBROUTINE SMOOTH computes the smooth array B corresponding to the unsmoothed array A. Both arrays contain N points. M, which must be odd, is the number of points in the running mean. In this smoothing routine, the beginning and the end points are not changed.

C.3.5 SUBROUTINE CO

At the altitude corresponding to the index j, SUBROUTINE CO computes the total ion density

$$\rho = \sum_{i=1}^{20} n_i m_i \quad (C14)$$

assuming the concentration is in diffusive equilibrium above the upper boundary.
That is,

$$\frac{dp}{dz} = - \rho g \quad (C6)$$

or

$$\frac{d(nkT)}{dz} = - nmg \quad (C7)$$

which reduces to

$$\frac{1}{n} \frac{dn}{dz} + \frac{1}{T} \frac{dT}{dz} = - \frac{mg}{kT} \quad (C8)$$

It is also assumed that above 400 km

$$\frac{dT}{dz} = 0.$$

Therefore, Eq. (C8) becomes

$$\frac{1}{n} \frac{dn}{dz} = - \frac{mg}{kT} \quad (C9)$$

An approximation to the acceleration of gravity at any altitude Z can be written as

$$g = \frac{g_0}{\left(1 + \frac{z}{R}\right)^2} = \frac{g_0 R^2}{(R + z)^2} \quad (C10)$$

where g_0 is the sea level value of the acceleration of gravity and R is the radius of the earth. Using Eq. (C10) and the assumption that the temperature is constant above the upper boundary, the solution to Eq. (C9) is

$$\ln\left(\frac{n}{n_0}\right) = - \frac{mg_0 R}{kT} \int_{z_0}^z \frac{dz}{(R + z)^2} \quad (C11)$$

have been computed for the species in the transport mode. It then calls SUBROUTINE CHEMION to compute the concentration profiles at the end of the current time step of all the remaining species. After all the new concentrations have been computed, the code then calls SUBROUTINE EFORU to compute the average velocities for those species in the transport mode.

Following this, a check is made to see if the solution for any species has converged at all altitudes. That is, is its largest relative change less than ERSPECI. If so, its IWHICH switch is set to zero and its ITER counter set to the current iteration number. The equations for this species will be skipped on the next iteration of the equation set.

A check is then made to see if the largest relative change among all the species is less than ERRMAX. If so, the iteration loop is exited and the solution accepted for this time step. If not, transfer is made back to the beginning of the DO 130 loop for the next iteration.

When the solution is accepted, the code then asks whether or not this solution should be sent to the output file. If printing has not been requested (PRINT=.FALSE.), transfer is made to that section of the code where the time step is doubled if possible, and the new concentrations and velocities are saved in extended core storage. Transfer is then made to STATEMENT 1000, where a new solution is generated over the next time step. If the solution is to be sent to the output file, the code then calculates all the subsidiary parameters to be printed out along with the concentrations and the average velocities. The code at this point is fairly explicit as to which quantities it is computing and sending to the output file. When all the information has been sent to the output file, the code then writes an interrupt file on Tape 4. This file contains all the information necessary to restart the code at this point at some later time. Since the development of a complete time dependent solution takes many hours of CDC 6600 CP time, it has to be accomplished in many short runs (about 4 hours each). The code is restarted from Tape 4 each time.

6.3.3 SUBROUTINE SMOOTHY

SMOOTHY controls the smoothing of those profiles that PROGRAM INTEG has determined should be smoothed. If no profiles are to be smoothed (NUM=0), then the subroutine is exited immediately upon entry. Because of the nature of the smoothing process, it is necessary to extrapolate the concentration profile above the upper boundary. The altitude above the upper boundary to which the extrapolation is done should be far enough away so that the effects of the smoothing at the top of the array are not felt at the upper boundary. In this code, the profiles are extrapolated to 500 km in 1=km=height steps. The extrapolation is accomplished

solution. This may be a new solution over the next time step or another try using a smaller time step because an error was encountered.

The code then saves the time, the solar zenith angle, the day-night transition height, and the day and night logical switches at the beginning of the time step. If an error is encountered later, these parameters will have to be reinitialized. The code then calls SUBROUTINE TIMER to compute the time and solar zenith angle at the end of the current time step. It then sets the JLOOK switch depending upon whether it is day or night. Following this, the IWHICH switches are all initialized to unity and the ITER locations to 999. Then, since the equations for helium will not be solved in this calculation, its IWHICH switch is set to zero.

The statement DO 130 is the beginning of the iteration loop. The loop beginning with the next statement is taken over all the species in the transport mode. The equations will be solved over the altitude extent for each species in succession. On the first pass through the loop, the quantities

$$\frac{1}{n_e} \frac{\partial n_e}{\partial Z} \approx \frac{1}{n_e} \frac{\Delta n_e}{\Delta Z} = \frac{n_{ej} - n_{ej-1}}{n_{ej} \Delta Z_{ej-1}} \quad (C5)$$

are computed and saved in the working array H. Next, the altitude dependent rate coefficients are computed using the current atmosphere and the values saved in extended core storage. If the JLOOK switch is on .TRUE., then the photodissociation rate coefficients are sent to the output file. If any of the concentration profiles need smoothing, this is done next and an acknowledgement is sent to the output file. To remain internally consistent, it is necessary now to recompute the velocity profiles for those species whose concentration profiles have been modified by smoothing. All the concentrations and velocities are then saved in extended core storage.

The code then brings into memory the values of the concentrations and velocities for the current species at the beginning of the time step. It then proceeds to compute a new concentration profile in the DO 100 loop. It also saves the largest relative change for this species in the BIGEST region. The relative change is defined as

$$\frac{n_{ej+1}^x - n_{ej}^x}{n_{ej+1}^x}.$$

The code then loops back to solve the continuity equation for the next species in turn. When the code exits at 105 CONTINUE, all the new concentration profiles

NSPEC is an array containing the order number from 1 to 56 of each species whose concentration is to be smoothed before the iteration procedure is started. The variable NUM indicates how many profiles are to be smoothed. If NUM=0, no smoothing will be done.

Two new labeled COMMONs are defined in PROGRAM INTEG.

C3.1 Labeled COMMON/LOGIC/

All the parameters listed in Labeled COMMON/LOGIC/ are logical variables. NIGHT is .TRUE. whenever the entire altitude region between the upper and lower boundaries is in darkness. Otherwise, it is .FALSE..

DAY is .TRUE. whenever the entire altitude region between the upper and lower boundaries is in daylight. Otherwise, it is .FALSE..

Both NIGHT and DAY variables are .FALSE. when the sun is rising or setting and only part of the altitude region between the upper and lower boundaries is in darkness.

JLOOK is set to .TRUE. to send to the output file, before the iteration begins, a table of the altitude distribution of the O_2 concentrations, the O_2 column densities, the O_3 concentrations, the O_3 column densities, and the rate coefficients for all the photodissociation reactions. If JLOOK is .FALSE., this table is not printed.

RESET is set to .TRUE. in SUBROUTINE TIMER whenever the time step has been recomputed because the next solution will be sent to the output file and the interrupt file.

END is set to .TRUE. whenever the computations for the current run are to be terminated.

C3.2 Labeled COMMON/COLUM/

The Labeled COMMON/COLUM/ contains information pertaining to the photodissociation rate coefficients.

02COL is used to transfer the O_2 column density from SUBROUTINE DISSOC to SUBROUTINE RATECN.

03COL is used to transfer the O_3 column density from SUBROUTINE DISSOC to SUBROUTINE RATECN.

DIS is the array where the photodissociation information is stored in SUBROUTINE RATECN for printing in PROGRAM INTEG. To avoid excessively long printouts, only every twelfth altitude value between the lower and the upper boundaries is stored in these arrays.

PROGRAM INTEG starts by initializing some of its variables. STATEMENT 1000 CONTINUE marks the beginning of the working code. After every successful or unsuccessful solution, the code will loop back to this statement to begin the next

and (9). This overlay uses two DIMENSION statements that contain the following variables:

PHI is an array EQUIVALENCED to the array U(1, 21) and is used as temporary storage of the total flux ϕ of each species during the printout phase of the code. The array U(1, 21) would normally contain the velocity profile of helium, but, since this species is held invariant in these calculations, these memory locations are available for other uses.

DED is an array EQUIVALENCED to the array E in COMMON. It is used as temporary storage of the turbulent diffusion velocity of each species successively during the printout phase of the code.

DEDZ is an array EQUIVALENCED to the array NZERO and is used as temporary storage of the vertical gradient of the turbulent velocity for each species during the printout phase of the code.

VBAR is an array EQUIVALENCED to the array F and is used as temporary storage of the molecular diffusion velocity of each species during the printout phase of the code. The molecular diffusion velocity of a species is equal to its average velocity minus the mean mass velocity.

H is an array used in various places in the code for temporary storage of arrays.

OUTP contains four words each of which contains one hollerith character of the local time in hours. It is computed by SUBROUTINE HOUR.

IWHICH is an array of integer switches, one for each species in the transport mode. If a switch is set to 1, the equations for that species will be iterated. If the switch is set to 0, the equations for that species will not be iterated. If IWHICH(21) is set to 0 before entering the iteration loop, the concentration of helium will remain fixed at its initial value throughout the calculations since its equations will not be iterated.

ITER is an array that stores the number of iterations required by each of the sets of equations before a final solution was obtained for a given time step.

BIGEST is an array that contains the largest relative difference in each species concentration profile. It contains the maximum value of

$$\frac{n_j^{z+1} - n_j^z}{n_j^z}$$

in the array as j varies from the lower boundary to the upper boundary. This array is recomputed after each iteration.

given time. The first file, which is slightly different from all the others, must be prepared before this code is first run. It contains the initial conditions on all the dependent variables. (See the discussion on the initial conditions in Section 7.1.) Since this is an unformatted file, it must be written exactly as it is read here. First, it reads the date that the first file was written, the file number (1 in this case), and the height in kilometers of the turbopause. It then skips over six words in the file that are not necessary to the current calculations. Then it reads the number of altitude steps in the complete arrays ($K = 692$), the number of altitude steps in the reduced arrays ($K_2 = 424$), and the altitude array in kilometers. After this, it reads the height step array, the fixed temperature array, the array $1/T \frac{dT}{dZ}$, and the fixed turbulent diffusion coefficient array. Next, it reads the initial velocities and initial concentrations.

In the following loop, the remainder of the interrupt files are skipped over and the tape is positioned at the beginning of a new file. At the completion of this loop, the locations containing the dependent variables will have been updated with the values obtained after the last successful solution. These values are then used as initial conditions for the next time step.

All files after file 1 on the interrupt tape are written in OVERLAY (2,0). Care must be taken not to write so many files on Tape 4 that it becomes full. Before this happens, a new Tape 4 must be made containing file 1 and the last file from the old Tape 4.

After the dependent variables have been initialized, the code initializes the solar zenith angle to the current time and computes the noontime solar zenith angle. It then reads from input Tape 5 and, where required, converts to proper units the wavelengths, solar fluxes, and absorption and ionization cross sections to be used in the photoionization calculations. Finally, from this same file, it reads the rate coefficient arrays for the energetic electron reactions. All that remains now is the calculation of the sunset (sunrise) solar zenith angles as a function of altitude. Also calculated are the times of sunset at the lower boundary and of sunrise at the upper boundary.

When all this initialization is complete, the concentration profiles corresponding to the initialized time are written to the output file. Finally, the current value of the variable TIMOUT is written to the output file in case the frequency of the output of solutions is to be changed for the next run.

C3. OVERLAY (2,0) PROGRAM INTEG

OVERLAY (2,0) contains all the programs that actually solve Eqs. (1), (7),

RADIUS (GRAV, ϕ)

$$= \frac{2 \cdot \text{GRAV}}{2.27 \times 10^{-9} \cos(2\phi) - 2 \times 10^{-2} \cos(4\phi) + 3.085426 \times 10^{-6}} \quad (\text{C4})$$

Next, the constants in labeled COMMON/CONSTAN/ are defined. After this, the constants for the particular solution to be generated are defined.

CLAT is the latitude in degrees.

DECL is the solar declination for the time of year when the calculations are to be made.

DIP is the magnetic dip angle appropriate to CLAT.

Next, the acceleration of gravity, the radius of the earth, and the constants required for later computing the solar zenith angles are computed. Information pertaining to the current calculations is then written to the output file (Tape 6).

The next section of this OVERLAY sets the chemical symbols along with the atomic or molecular weight, the mixing ratio, the collision cross section, the polarizability, and the thermal diffusion factor for the 26 transport species. Following this, the symbols for the species computed from chemistry only are written. The atomic or molecular weight of each species is then converted to mass in grams. The ground level mixing ratio of N₂ is computed to ensure that the sum of all the mixing ratios is identically unity. The mixing ratios of the species are not necessary for the code described here. However, if other initial conditions are to be computed as discussed in Keneshea et al,⁸ then the values of the mixing ratios are required.

The following section of INITIAL obtains input information from various sources. The list of chemical reactions with their rate coefficients must be prepared separately and saved for use here as the file called Tape 5. This information is coded according to the FORMAT defined in statements 770 or 771, depending upon whether the reaction is not height dependent (770) or is height dependent (771). As the reactions are read in, they are also sent to the OUTPUT file (Tape 6). After the reaction set is read in, the code reads in the photodissociation rate coefficient tables (Tables B4 through B20). All values in these tables are in log base 10. These tables also are computed and saved separately. Since these numbers are functions only of the O₂ and O₃ column densities, they are universal. They need be computed only once and can then be input to any calculation regardless of geographic location, season of the year, or time of day.

After the photodissociation rate coefficients are initialized, the invariant part of the collision frequencies ν_{ij} are computed and saved in the D array.

The next section of the code reads the interrupt file Tape 4. This tape is structured in files, each containing the solution for all the dependent variables at a

morning or evening twilight periods and the current altitude is below the day-night transition altitude, then the computation of the rate coefficients is skipped. Otherwise, they are computed.

During the daytime, SUBROUTINE DISSOC is called for the computation of the photodissociation rate coefficients. These are then stored in the proper cells of the DK array. If the JLOOK switch is on .TRUE., these rate coefficients along with the concentrations and column densities of O_2 and O_3 are stored in the DIS region for later printing in the output file. Following this, the rate coefficients for the energetic electron reactions are computed. The values in the tables are interpolated to the current solar zenith angle and altitude. Next, the column densities of O_2 , N_2 , and O at the current solar zenith angle are approximated. These column densities are then used to compute the EUV attenuated flux at each wavelength. The rate coefficients for the photoionization processes are then computed. All EUV radiation below 0.15 nm is considered to be X radiation, which is handled differently from the longer EUV radiation. The total x-ray ionization is computed and partitioned among the species according to Swider.¹⁹ Finally, the total EUV rate coefficients are computed as well as the rate coefficients for the ionization of nitric oxide by Lyman alpha and the ionization of $O_2(^1\Delta g)$.

If the code is in the nighttime, the above computations are skipped and the code continues from this point. The contributions of the scattered radiations to the ionization rates are computed next. Finally, the contribution of cosmic rays is added into the rate coefficients and the subroutine is exited.

3.15 FUNCTION YINT

FUNCTION YINT performs interpolation where it is required. This function essentially computes a cubic spline fit to the parameter under consideration. It then returns the interpolated value in YINT corresponding to XII in the (X, Y) arrays.

3.16 SUBROUTINE CHAPMAN

SUBROUTINE CHAPMAN computes an approximation to the Chapman function, which is required in computing the attenuation of solar energy in its transmission through the atmosphere. The approximations computed are those given by Swider and Gardner.^{C2}

C2. Swider, W. Jr., and Gardner, M.E. (1967) On the Accuracy of Certain Approximations for the Chapman Function, AFCRL-67-0468, AD 658826.

C3.17 SUBROUTINE DISSOC

SUBROUTINE DISSOC returns the value of the photodissociation rate coefficients appropriate to the current solar zenith angle and the altitude whose index is IALT. The rate coefficients for all species except nitric oxide are determined as follows: The column densities of O_2 and O_3 are first approximated for the current O_2 and O_3 concentration distributions in the code. A linear interpolation is then made in the appropriate photodissociation table to determine the current photodissociation rate coefficient for each of the photodissociation processes. Finally, using the data of Cieslik and Nicolet^{C3} as given in the DATA statement of this subroutine, the photodissociation rate coefficient for nitric oxide is computed.

C3.18 FUNCTION TRPLT2

FUNCTION TRPLT2 interpolates in the two-dimensional photodissociation rate coefficient tables. The two-dimensional array to be interpolated is fed to this function through the parameter ARRAY. The dimensions of ARRAY are (N, M). The indices of elements in ARRAY are IO2 and IO3. The increments required to perform the interpolation are DELO2 and DELO3.

C3.19 FUNCTION TRPLT1

FUNCTION TRPLT1 interpolates in the one-dimensional photodissociation rate coefficient tables. The array to be interpolated is fed to the function through the parameter ARRAY. The index in ARRAY is IO3 and the interpolation increment is DELO3.

C3.20 SUBROUTINE INDEX

SUBROUTINE INDEX computes an index I and an increment DEL corresponding to the column density COL. LIM is the lowest power of 10 of the column densities of O_2 and O_3 in the photodissociation tables under consideration.

C3.21 SUBROUTINE CHEMPR

Each time SUBROUTINE CHEMPR is called, it sends to the output file the complete list of chemical reactions for the altitude whose index is IZ. It also sends the rate coefficient for each reaction and its forward rate. This subroutine is also written by the WRITER codes given in Keneshea.³ The output from this subroutine is used to determine the relative importance of the chemical reactions in different altitude regions.

^{C3} Cieslik, S., and Nicolet, M. (1973) The aeronomic dissociation of nitric oxide, *Planet. Space Sci.* 21:925-938.

C3.22 SUBROUTINE HOUR

SUBROUTINE HOUR converts the value of TIME, which is in accumulated seconds from the noontime of the day on which the simulation was started, to the local time in hours (0000 to 2359). The digits of the hour are returned in hollerith in successive words of the OUTP array.

```

C-EPLA-1992-1
PROGRAM MAIN. INPUT=60, OUTPUT=51, TAPES=55, TAPE5=56.

      COMMON/ N(692,30), CONC(424,26), U(692,26), NZERO(692), JZERO(692)
      COMMON/ FORM(692), REMV(692), E(692), F(692), Z(692),
      2      DEDY(692), TEMP(692), DENDZ(692), DELTAZ(692), SUNSET(692)
      COMMON/ GRAN, COJO, SIND, TIME, CX1, RADIUS, ISPECI, KSPECI, NRAC,
      ;      K2, KM1N, ITUPB, CXINDN, SCY1, NRAC2
      COMMON/ ESM/ C(692,30), OLUCJN(424,26), OLUU(692,26), RATES(88,692)
      COMMON/ TIMES, DAYS, DELT, TMAX, TSET, TRISE
      COMMON/ AL1110/, ZBOT, ZTOP, TURB, ZTRANS
      COMMON/ RESTART, IFRAME, TOFF, TYME, TODAY
      COMMON/ OUTPUTS/ TIMEOUT, OPRINT, PRINT
      COMMON/ ERRORS/ ERMAX, ERSPEC1, ITMAX
      COMMON/ CCUSTAN, PI, RADDEG, DEGRAD, RADSEC, BOLTZ, SIN2I, ATCON
      COMMON/ JAYS/ J192(26,25), J193(26,26), J194(22), J195(32,28),
      1      J196(31,30), J197(27,28), J198(26,28), J200(19,15),
      2      J201(26,27), J202(22), J203(30,29), J204(35,28), J205(22)
      COMMON/ COLFRE/ D(26,26)
      COMMON/ ATCON, MASS/ ATWT(26), THERM(26), SYMBOL(56), POLAR(26)
      COMMON/ RATECON, AK/ (88), BK(88), CK(83), DK(215), CYMB(6,215)
      COMMON/ IONS/ MAYE(88), FLUX(88), SIGMA(9,88), R(8,4,8)
      REAL N, NZERO
      LEVEL 3, CLOX, OLDCON, OLUU, RATES

      !SPECI IS THE TOTAL NUMBER OF SPECIES IN THE MODEL. KSPECI IS
      !THE NUMBER OF SPECIES IN THE TRANSPORT MODE WITH ALL THE
      !OTHERS IN THE CHEMISTRY MODE.
      LSPECI=56
      KSPECI=26
      NRAC=127
      NRAC2=68

      C   ALITUDES
      C   LSPECI=56
      C   ZTOP=100,
      C   ZB=12,
      C   ZL=12,
      C   ZD=42;

```

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      C   MIONEUT
      C   MIONEUT 59
      C   MIONFUT 60
      C   MIONFUT 61
      C   MIONEUT 62
      C   MIONEUT 63
      C   MIONEUT 64
      C   MIONEUT 65
      C   MIONEUT 66
      C   MIONEUT 67
      C   MIONEUT 68
      C   MIONEUT 69
      C   MIONEUT 70
      C   MIONEUT 71
      C   MIONEUT 72
      C   MIONEUT 73
      C   MIONEUT 74
      C   MIONEUT 75
      C   MIONEUT 76
      C   MIONEUT 77
      C   MIONEUT 78
      C   MIONEUT 79
      C   MIONEUT 80
      C   MIONEUT 81
      C   MIONEUT 82
      C   MIONEUT 83
      C   MIONEUT 84
      C   MIONEUT 85
      C   MIONEUT 86
      C   MIONEUT 87
      C   MIONEUT 88
      C   MIONEUT 89
      C   MIONEUT 90
      C   MIONEUT 91
      C   MIONEUT 92
      C   MIONEUT 93
      C   MIONEUT 94
      C   MIONEUT 95
      C   MIONEUT 96
      C   MIONEUT 97
      C   MIONEUT 98
      C   MIONEUT 99
      C   MIONEUT 100
      C   MIONEUT 101
      C   MIONEUT 102
      C   MIONEUT 103
      C   MIONEUT 104
      C   MIONEUT 105
      C   MIONEUT 106
      C   MIONEUT 107
      C   MIONEUT 108

C   RESTART
C   TOFF=250.
C   IFRAME=1
C   TIME=14400.
C   CALL DATE(TODAY)
C   WRITE(6,50) TODAY
C   TIMES
C   TIME=0.
C   TIMEOUT=0.
C   DELT=2.0**(-10)
C   TMAX=1800.
C   DAYS=40.
C   OUTPUTS
C   OPRINT=12.
C   ERRORS
C   ERMAX=0.3
C   ERSPECI=0.01
C   ITMAX=100
C   CALL OVERLAY(6HIONEUT,1,0)
C   CALL OVERLAY(6HIONEUT,2,0)
C   START COMPUTING TIME-HEIGHT PROFILES.
C   CALL MOVLEV(N,OLDN,20760)
C   CALL MOVLEV(CONC,OLDCON,11024)
C   CALL MOVLEV(U,OLDU,17992)
C   REWIND 4
C   50 FORMAT(1H1,* THIS RUN WAS MADE ON *,A10)
C   END

```

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2      IIONEUT
3      IIONEUT
4      IIONEUT
5      IIONEUT
6      IIONEUT
7      IIONEUT
8      IIONEUT
9      IIONEUT
10     IIONEUT
11     IIONEUT
12     IIONEUT
13     IIONEUT
14     IIONEUT
15     IIONEUT
16     IIONEUT
17     IIONEUT
18     IIONEUT
19     IIONEUT
20     IIONEUT
21     IIONEUT
22     IIONEUT
23     IIONEUT
24     IIONEUT
25     IIONEUT
26     IIONEUT
27     IIONEUT
28     IIONEUT
29     IIONEUT
30     IIONEUT
31     IIONEUT
32     IIONEUT
33     IIONEUT
34     IIONEUT
35     IIONEUT
36     IIONEUT
37     IIONEUT
38     IIONEUT
39     IIONEUT
40     IIONEUT
41     IIONEUT
42     IIONEUT
43     IIONEUT
44     IIONEUT
45     IIONEUT
46     IIONEUT
47     IIONEUT
48     IIONEUT
49     IIONEUT
50     IIONEUT
51     IIONEUT
52     IIONEUT
53     IIONEUT
54     IIONEUT
55     IIONEUT
56     IIONEUT
57     IIONEUT

C  DIVERGENCE, J2
C  PROGRAM INIT.A
C  DIME ISIGN S(26),RMIX(26),SIG(88)
C
C  COMMON N(692, 30),CONC(424,26),U(692,26),NZERO(692),UZERO(692)
C
C  COMMON FORM(692),REMV(692),E(692),F(692),Z(692),
C  2      DEDY(592),TEMP(692),DTENDZ(692),DELTAZ(692),SUNSET(692)
C
C  COMMON GRAV,COSD,SIND,TIME,CX1,RADIUS,ISPECI,KSPECI,NREAC,K,
C  1      K2,KMIN1,ITURB,CXINDON,SCXI,NREAC
C
C  COMMON/ATCONS/AINT(26),MASS(26),THERM(26),SYMBOL(56),POLAR(26)
C
C  COMMON/TIMES/DAYS,DELT,TMAX,TSET,TRISE
C
C  COMMON/ALTITUO/ZBOT,ZTOP,TURB,ZTRANS
C
C  COMMON/RESTART/IFRAME,TOFF,TYME,TODAY
C
C  COMMON/CONSTAN/PI,RADDEG,DEGRAD,RADSEC,BOLTZ,SIN21,ATCON
C
C  COMMON/JAYS/J192(26,25),J193(28,26),J194(22),J195(32,28),
C  1      J196(31,30),J197(27,28),J198(26,28),J200(19,15),
C  2      J201(26,27),J202(22),J203(30,29),J204(35,28),J205(22)
C
C  COMMON/IONS/WAVE(88),FLUX(88),SIGMA(9,88),R(8,4,8)
C
C  COMMON/RATECON/AK(88),BK(88),CK(88),DK(215),CYMB(6,215)
C
C  COMMON/COLFRE/D(26,26)
C
C  REAL N,MASS,J192,J193,J194,J195,J196,J197,J198,J200,J201,J202
C  REAL J203,J204,J205
C
C  THE FOLLOWING DATA STATEMENT CONTAINS THE PHOTOIONIZATION
C  CROSS SECTIONS FOR ATOMIC NITROGEN SCALED BY 1E18.
C  SEE "AERONOMY" BY BANKS AND KOCHARTS, PART A, P102-104.
C
C  DATA SIG/21*(0.),20*(10.0),2*(10.5),2*(10.4),10.75,4*(11.0),
C  1      11.8,11.9,2*(12.0),11.9,11.5,11.1,11.0,9.3,10.0,9.7,
C  2      8.5,8.8,2*(8.0),7.3,6.7,6.0,5.3,6.3,4.3,3.4,2.8,2.2,1.3,
C  3      0.58,0.45,0.36,0.32,0.23,0.16,0.093,0.04,0.03,0.015,
C  4      0.006,0.001,0.0001/
C
C  FUNCTION DEFINITION FOR SEA LEVEL ACCELERATION OF GRAVITY AS
C  A FUNCTION OF LATITUDE IN CM/SEC**2.
C
C  GRAVITY(ALAT)=978.0356*(1.0+0.0052885*(SIN(ALAT))**2
C  1      -5.9E-06*(SIN(2.0*ALAT))**2)
C
C  FUNCTION DEFINITION FOR RADIUS OF THE EARTH IN KILOMETERS AS
C  A FUNCTION OF LATITUDE AND SEA LEVEL GRAVITY.
C
C  C

```

```

      RADUS(GRAV,ALAT)*(2.0*GRAV*1.0E-5)/(2.27E-09*COS(2.0*ALAT))  

      -2.0E-12*COS(4.0*ALAT)+3.085462E-06) 59  

      LIONEUT 60  

      LIONEUT 61  

      LIONEUT 62  

      LIONEUT 63  

      LIONEUT 64  

      LIONEUT 65  

      LIONEUT 66  

      LIONEUT 67  

      LIONEUT 68  

      LIONEUT 69  

      LIONEUT 70  

      LIONEUT 71  

      LIONEUT 72  

      LIONEUT 73  

      LIONEUT 74  

      LIONEUT 75  

      LIONEUT 76  

      LIONEUT 77  

      LIONEUT 78  

      LIONEUT 79  

      LIONEUT 80  

      LIONEUT 81  

      LIONEUT 82  

      LIONEUT 83  

      LIONEUT 84  

      LIONEUT 85  

      LIONEUT 86  

      LIONEUT 87  

      LIONEUT 88  

      LIONEUT 89  

      LIONEUT 90  

      LIONEUT 91  

      LIONEUT 92  

      LIONEUT 93  

      LIONEUT 94  

      LIONEUT 95  

      LIONEUT 96  

      LIONEUT 97  

      LIONEUT 98  

      LIONEUT 99  

      LIONEUT 100  

      LIONEUT 101  

      LIONEUT 102  

      LIONEUT 103  

      LIONEUT 104  

      LIONEUT 105  

      LIONEUT 106  

      LIONEUT 107  

      LIONEUT 108  

      LIONEUT 109  

      LIONEUT 110  

      LIONEUT 111  

      LIONEUT 112  

      LIONEUT 113  

      LIONEUT 114  

      LIONEUT 115  

      C CONSTANTS  

      C  

      C PI=3.14159265359  

      RADDEG=PI/180.  

      DEGRAD=180./PI  

      RADSEC=PI/43200.  

      BOLTZ=1.38047E-16 65  

      C  

      C SET CONSTANTS FOR PARTICULAR MODEL CALCULATION.  

      C  

      CLAT=30.0  

      DECL=-19.55  

      DIP=60.0  

      PHILAT=CLAT+RADDEG  

      GRAV=GRAVITY(PHILAT)  

      RADIUS=RADIUS(GRAV,PHILAT)  

      DECLAT=DECL+RADDEG  

      COSD=COS(DECLAT)*COS(PHILAT)  

      SIND=SIN(DECLAT)*SIN(PHILAT)  

      SIN2I=(SIN(DIP+RADDEG))**2  

      WRITE(6,25) CLAT,DECL,DIP,GRAV,RADIUS  

      25 FORMAT(1HO,5X,"LATITUDE OF MODEL **,OPF6.2.", DEGREES. /  

      16X,"SOLAR DECLINATION **,F6.2.", DEGREES. /6X,"MAGNETIC DIP **/  

      2*ANGLE **,F6.2., DEGREES. /6X,"ACCELERATION OF GRAVITY **,1PE12.5.  

      3* CM/SEC**2. /6X,"RADIUS OF THE EARTH **,E12.5., KM.")  

      C  

      C SET ATOMIC AND MOLECULAR WEIGHTS, MIXING RATIOS, COLLISION CROSS  

      C SECTIONS, THERMAL DIFFUSION FACTORS AND POLARIZABILITIES.  

      C  

      ATCON=1.66035E-24  

      DO 60 I=1,KSPECI  

      THERM(I)=1.0  

      POLAR(I)=0.0  

      RMIX(I)=0.0  

      60 CONTINUE  

      SYMBOL(1)=10H H+  

      ATWT(1)=1.00797  

      S(1)=3.04E-08  

      SYMBOL(2)=10H HE+  

      ATWT(2)=4.0026  

      S(2)=2.04E-08  

      SYMBOL(3)=10H O+  

      ATWT(3)=15.9994  

      S(3)=1.2E-08  

      SYMBOL(4)=10H O2+  

      ATWT(4)=31.9988  

      S(4)=3.3E-08  

      SYMBOL(5)=10H NO+  

      ATWT(5)=30.006132  

      S(5)=3.4E-08  

      SYMBOL(6)=10H N+  

      ATWT(6)=14.026732  

      S(6)=1.42E-08  

      SYMBOL(7)=10H O 85  


```

```

116 1.0E-011
117 1.0E-011
118 1.0E-011
119 1.0E-011
120 1.0E-011
121 1.0E-011
122 1.0E-011
123 1.0E-011
124 1.0E-011
125 1.0E-011
126 1.0E-011
127 1.0E-011
128 1.0E-011
129 1.0E-011
130 1.0E-011
131 1.0E-011
132 1.0E-011
133 1.0E-011
134 1.0E-011
135 1.0E-011
136 1.0E-011
137 1.0E-011
138 1.0E-011
139 1.0E-011
140 1.0E-011
141 1.0E-011
142 1.0E-011
143 1.0E-011
144 1.0E-011
145 1.0E-011
146 1.0E-011
147 1.0E-011
148 1.0E-011
149 1.0E-011
150 1.0E-011
151 1.0E-011
152 1.0E-011
153 1.0E-011
154 1.0E-011
155 1.0E-011
156 1.0E-011
157 1.0E-011
158 1.0E-011
159 1.0E-011
160 1.0E-011
161 1.0E-011
162 1.0E-011
163 1.0E-011
164 1.0E-011
165 1.0E-011
166 1.0E-011
167 1.0E-011
168 1.0E-011
169 1.0E-011
170 1.0E-011
171 1.0E-011
172 1.0E-011

120
  RMX(7)=1.0E-08
  RMIX(7)=1.0E-08
  S(7)=2.65E-08
  SYMBOL(8)=10H 02
  THERM(8)=1.0+0.0
  POLAR(8)=1.59
  ATWT(8)=31.9988
  RMIX(8)=0.209545
  S(8)=3.40E-08
  SYMBOL(9)=10H 03
  ATWT(9)=47.9982
  RMIX(9)=7.0E-06
  S(9)=4.0E-08
  SYMBOL(10)=10H OH
  ATWT(10)=17.00737
  RMIX(10)=1.0E-10
  S(10)=3.0E-08
  SYMBOL(11)=10H H
  THERM(11)=1.0-0.39
  POLAR(11)=0.667
  ATWT(11)=1.00797
  RMIX(11)=1.6E-13
  S(11)=3.04E-08
  SYMBOL(12)=10H HD2
  ATWT(12)=33.00677
  RMIX(12)=1.0E-12
  S(12)=4.0E-08
  SYMBOL(13)=10H H2D
  ATWT(13)=18.01534
  RMIX(13)=5.0E-06
  S(13)=3.5E-08
  SYMBOL(14)=10H H2D2
  ATWT(14)=34.01474
  RMIX(14)=1.0E-10
  S(14)=4.0E-08
  SYMBOL(15)=10H H2
  THERM(15)=1.0-0.31
  POLAR(15)=0.82
  ATWT(15)=2.01594
  RMIX(15)=4.5E-07
  S(15)=2.97E-08
  SYMBOL(16)=10H N
  POLAR(16)=1.1
  ATWT(16)=14.006732
  RMIX(16)=1.0E-11
  S(16)=1.42E-08
  SYMBOL(17)=10H N2D
  POLAR(17)=1.1
  ATWT(17)=14.006732
  RMIX(17)=1.0E-16
  S(17)=1.42E-08
  SYMBOL(18)=10H NQ
  ATWT(18)=1.74
  RMIX(18)=30.00532
  S(18)=1.0E-11

```

```

      5,18)=3.4E-03  

SYMBOL(19)=1OH NO2  

ATWT(19)=-4E-035532  

RMIX(19)=1.0E-11  

S(19)=3.8E-08  

SYMBOL(20)=1OH C21D  

FOLAR(20)=1.59  

ATWT(20)=31.9963  

RMIX(20)=2.3E-C6  

S(20)=3.3E-08  

SYMBOL(21)=1OH HELIUM  

THERM(21)=1.0E-0.36  

POLAR(21)=0.21  

ATWT(21)=4.0026  

RMIX(21)=5.24E-06  

S(21)=2.04E-08  

SYMBOL(22)=1OH ARGON  

ATWT(22)=39.948  

RMIX(22)=0.00934  

S(22)=2.97E-08  

SYMBOL(23)=1OH N2  

POLAR(23)=1.76  

ATWT(23)=28.013464  

RMIX(23)=0.78  

S(23)=3.15E-08  

SYMBOL(24)=1OH N2O  

POLAR(24)=3.00  

ATWT(24)=4.012864  

RMIX(24)=2.7E-07  

S(24)=3.7E-08  

SYMBOL(25)=1OH CO2  

POLAR(25)=2.63  

ATWT(25)=4.00995  

RMIX(25)=3.14E-04  

S(25)=3.8E-08  

SYMBOL(26)=1OH CO  

POLAR(26)=1.97  

ATWT(26)=28.01055  

RMIX(26)=1.0E-08  

S(26)=3.0E-08

```

C THE FOLLOWING SPECIES ARE COMPUTED FROM CHEMISTRY ONLY.

C

173 IIONEUT
174 IIONEUT
175 IIONEUT
176 IIONEUT
177 IIONEUT
178 IIONEUT
179 IIONEUT
180 IIONEUT
181 IIONEUT
182 IIONEUT
183 IIONEUT
184 IIONEUT
185 IIONEUT
186 IIONEUT
187 IIONEUT
188 IIONEUT
189 IIONEUT
190 IIONEUT
191 IIONEUT
192 IIONEUT
193 IIONEUT
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      SYMBOL( 42 ) = 1H H3O+
      SYMBOL( 43 ) = 1H H7O3+
      SYMBOL( 44 ) = 1H H9O4+
      SYMBOL( 45 ) = 1H H11O5+
      SYMBOL( 46 ) = 1H H3O+ . N2
      SYMBOL( 47 ) = 1H H3O+ . OH
      SYMBOL( 48 ) = 1H H3O+ . CO2
      SYMBOL( 49 ) = 1H H5O2+ . CO2
      SYMBOL( 50 ) = 1H O2+ . H2O
      SYMBOL( 51 ) = 1H NO+ . H2O
      SYMBOL( 52 ) = 1H NO+ . 2H2O
      SYMBOL( 53 ) = 1H NO+ . 3H2O
      SYMBOL( 54 ) = 1H NO+ . CO2
      SYMBOL( 55 ) = 1H HNO2
      SYMBOL( 56 ) = 1H NO3
      SUMMIX=0.
      DO 2 I=1,KSPECI
      MASS(I)=ATWT(I)*ATCON
      IF(I .EQ. 23) GO TO 2
      SUMMIX=SUMMIX+RMIX(I)
      CONTINUE
      RMIX(23)=1.0-SUMMIX
      2
      C READ CHEMICAL REACTIONS AND RATE CONSTANTS.
      C
      245      WRITE(6,650)
      650      FORMAT(1H0, * THE CHEMICAL REACTIONS USED IN THIS RUN ARE::*)
      DO 765 J=1,NREAC
      READ(5,770) IREAC,(CYMB(I,J),I=1,6),DK(J)
      770      FORMAT(13.5A10,A5,1PE8.2)
      WRITE(6,771) IREAC,(CYMB(I,J),I=1,6),DK(J)
      771      FORMAT(2X,13.5A10,A5,1PE9.2)
      765      CONTINUE
      DO 35 J=1,NREAC2
      READ(5,772) IREAC,(CYMB(I,NREAC+J),I=1,6),AK(J),BK(J),CK(J)
      772      FORMAT(13.5A10,A5,1PE8.2,1PE9.2)
      WRITE(6,773) IREAC,(CYMB(I,NREAC+J),I=1,6),AK(J),BK(J),CK(J)
      773      FORMAT(2X,13.5A10,A5,1PE9.2,0PF6.2,1PE10.2)
      35      CONTINUE
      C READ PHOTODISOCIATION RATE COEFFICIENTS AS LOG BASE 10.
      C
      255      REMIND 3
      READ(3) ((J193(L,M),L=1,28),M=1,26)
      READ(3) ((J192(L,M),L=1,26),M=1,25)
      READ(3) ((J196(L,M),L=1,31),M=1,30)
      READ(3) ((J197(L,M),L=1,27),M=1,28)
      READ(3) ((J195(L,M),L=1,32),M=1,28)
      READ(3) ((J194(L),L=1,22))
      READ(3) ((J200(L,M),L=1,19),M=1,15)
      READ(3) ((J205(L),L=1,22))
      READ(3) ((J201(L,M),L=1,26),M=1,27)
      READ(3) ((N(L,M),L=1,26),M=1,25)
      READ(3) ((J202(L),L=1,22))
      READ(3) ((N(L,M),L=1,26),M=1,23)
      READ(3) ((N(J,M),L=1,30),M=1,30)
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      7285
      7290
      7295
      7300
      7305
      7310
      7315
      7320
      7325
      7330
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      7365
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      7385
      7390
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      7575
      7580
      7585
      7590
      7595
      7600
      7605
      7610
      7615
      7620
      7625
      7630
      
```

```

      READ(3) ((J203(L,M),L=1,30),M=1,29)
      READ(3) ((J204(L,M),L=1,35),M=1,28)
      READ(3) ('N(L,M),L=1,37),M=1,20)
      READ(3) ((J198(L,M),L=1,26),M=1,28)

290   C COMPUTE INVARIANT PART OF COLLISION FREQUENCIES.
      C
      DO 22 I=1,KSPECI
      DO 22 J=1,KSPECI
      READ(4) DATEOF,IFRAME,TURB,(N(J,1),J=1,6),K,K2,(Z(J),J=1,K)
      D(J,1)=0.5*((S(I)+S(J))*2)*SQRT((2.0*PI*BOLTZ*(MASS(I)+MASS(J)))
      1          /(MASS(I)*MASS(J))))
      22  CONTINUE
      IFRAM=IFRAME-1

      REWIND 4

300
      READ(4) DATEOF,IFRAME,TURB,(N(J,1),J=1,6),K,K2,(Z(J),J=1,K)
      READ(4) (DELTAZ(L),L=1,K)
      READ(4) (TEMP(L),L=1,K)
      READ(4) (ITEMOD(L),L=1,K)
      READ(4) (DTEMOD(L),L=1,K)
      READ(4) (DEDY(L),L=1,K)
      DO 1 J=1,K
      DEDY(J)=3.0*DEDY(J)
      1  CONTINUE
      DO 5 L=1,KSPECI
      READ(4) (U(J,L),J=1,K)
      5  CONTINUE
      DO 20 L=1,ISPECI
      IF(L.GT.30) GO TO 10
      READ(4) (N(J,L),J=1,K)
      GO TO 20
      READ(4) (COND(J,L=30),J=1,K2)
      20  CONTINUE
      KMIN1=K-1
      13  READ(4)
      IF.EOF(4) 14,13
      IF(IFRAM.EQ.0) GO TO 15
      DO 140 I=1,IFRAM
      READ(4) DATEOF,IFRAME,TIME,SCXI,CXI,DELT,VINOUT
      DO 18 L=1,KSPECI
      READ(4) (U(J,L),J=1,K)
      CONTINUE
      DO 19 L=1,ISPECI
      IF(L.GT.30) GO TO 11
      READ(4) (N(J,L),J=1,K)
      GO TO 19
      READ(4) (COND(J,L=30),J=1,K2)
      19  CONTINUE
      16  READ(4)
      IF.EOF(4) 140,16
      140 CONTINUE
      15  CXI=ACOS(COSD+COS(RADSEC+TIME)+SIND)
      CXIND=ACOS(COSD+SIND)
      SCXI=CXI*DEGRAD
      DO 4 J=1,K
      IF(Z(J).GT.TURB) GO TO 6
      4  CONTINUE
      6  ITURB=J-1
      C

```

```

C READ WAVELENGTH, L, L AND CROSS SECTION DATA FOR PHOTOIONIZATION.
C
345      READ(5,200) L,WAVE(L),FLUX(L),(SIGMA(I,L),I=1,6),J=1,88)
        FORMATTAB,3+L,4
        READ(5,205) J,FLUX(J),J=84,88)
144      IONNEUT
145      IONNEUT
146      IONNEUT
147      IONNEUT
148      IONNEUT
149      IONNEUT
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400      IONNEUT

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400      N1=12
        N2=22
        CONTINUE
        WRITE(6,605) TIME,SCX1,IFRAME
        WRITE(6,315) SYMBOL(J),J=23,30)
        WRITE(6,611) (SYMBOL(J),J=23,30)
        WRITE(6,320) (Z(J),(N(J,I),I=23,30),TEMP(J),DTENDZ(J),DEDY(J),
        1J=1,K,12)
        WRITE(6,320) Z(KMIN1),N(KMIN1),I=23,30),TEMP(KMIN1).
        DTENDZ(KMIN1),DEDY(KMIN1)
        WRITE(6,320) Z(K),(N(K,I),I=23,30),TEMP(K),DTENDZ(K),DEDY(K)
        FORMAT(*,ALT(KM)*,4X,8A10,*TEMP *(1/T)DT/DZ K-EDDY*)
        WRITE(6,605)
        WRITE(6,315) TIME,SCX1,IFRAME
        WRITE(6,125) SYMBOL(M),M=31,43)
        FORMAT(*,ALT(KM)*,13A9)
        WRITE(6,130) (Z(J),(CONC((J,M),M=1,13),J=1,K2,12)
        WRITE(6,130) Z(K2),(CONC(K2,M),M=1,13)
        FORMAT(OPFB-2,1X,1P13E9.2)
        WRITE(6,605)
        WRITE(6,315) TIME,SCX1,IFRAME
        WRITE(6,125) (SYMBOL(M),M=44,56)
        WRITE(6,130) (Z(J),(CONC((J,M),M=14,26),J=1,K2,12)
        WRITE(6,130) Z(K2),(CONC(K2,M),M=14,26)
        FORMAT(1H1,60X17HINITIAL PROFILES.)
        FORMAT(*,ALT(KM)*,11A10)
        315  FORMAT(1H ,10X,*TIME **,1PE12.5,* SEC.**,10X,*CX1 **,E12.5,
        1* DEGREES,**,10X,*FRAME NO.**,I4)
        320  FORMAT(OPFB-2,2X,1P11E10.3)
        WRITE(6,325) TIMEOUT
        325  FORMAT(1H0,* TIMEOUT **,1PE13.5)
        END

```

```

OVERLAY(2,0)
PROGRAM IN,EL
C
C      DIMENSION PHIN(1),DFOZ(1),DFDZ(1),VBARI(1),H(692),
C      DIMENSION QUTR(4),IWHICH(26),ITER(26),BIGEST(26),NSPEC(56)
C
C      COMMON/N692/DO1,CNC(424,26),U(692,26),NZERD1S92),UZERO(692)
C
C      COMMON FGRM(692),REMV(692),E(692),F(692),Z(692),
C      2      DDAY(692),TEMP(692),DTENDZ(692),DETAZ(692),SUNSET(692)
C
C      COMMON GRAV,CLSD,SIND,TIME,CXI,RADIUS,ISPEC1,KSPEC2,NRECAC,K,
C      1      K2,MINI,ITURB,CXINUN,SCX1,NRECAC2
C
C      COMMON/ESXN/UDCN(692,30),CLDCON(424,26),OLDU(692,26),RATES(88,692)
C
C      COMMON/ATMFLG/ DAYS,DELT,TMAX,TSET,TRISE
C
C      COMMON/ACTFLG/ZSC,ZNOP,TURB,ZTRANS
C
C      COMMON/RESTART/IFRAME,TOFF,TYME,TODAY
C
C      COMMON/OUTPUTS/TIMOUT,OPRINT,PRINT
C
C      COMMON/ERRORS/ERRMAX,ERSPEC1,ITMAX
C
C      COMMON/DIGIC/NIGHT,DAY,JLOOK,RESET,END
C
C      COMMON/ATCONS/ATW(26),MASS(26),TERM(26),SYMBOL(56),POLAR(26)
C
C      COMMON/CONSTANT/PI,RADdeg,DEGRAD,RADSEC,BOLTZ,SINC1,ATCIN
C
C      COMMON/COLONY/D2CDL,03CC1,DIS(15,59)
C
C      COMMON/RATECON/AK(88),BK1S81,CK(88),DK(215),CYMB(6,215)
C
C      LEVEL 3, OLD, OLDCN,OLDU, RATES
C      REAL N,NZERO,NEWN,MASS,MBAR,NTOTAL
C      LOGICAL JLOOK,NIGHT,DAY,RESET,PRINT,END,LASTNT,LASTDA
C      EQUIVALENCE (U(1,21) H,PHI),(E,DED),(F,VBAR),(NZERO,DEDZ)
C      EQUIVALENCE (UZERO,DFDZ)
C
C      DATA NSETCH/1/
C
C      FUNCTION DEFINITION FOR CORRECTION TO ACCELERATION OF GRAVITY.
C
C      GRAYCOR(ZZ)=C/(1.0+ZZ/RADIUS)**21
C
C      IF ANY SMOOTHING IS TO BE DONE, SET NUM TO THE NUMBER OF SPECIES
C      TO BE SMOOTHED AND PUT THEIR SPECIFIC NUMBERS INTO THE NSPEC REGION.
C      IF NO SMOOTHING IS TO BE DONE, SET NUM TO ZERO.
C
C      NUM=0
C      ENDIF, FALSE.
C      NCHECK=1
C      DO1=1,26,1
C      SCX1=1
C      NZERD1S92=1
C      NZERD1=1
C      NZERD2=1
C      NZERD3=1
C      NZERD4=1
C      NZERD5=1
C      NZERD6=1
C      NZERD7=1
C      NZERD8=1
C      NZERD9=1
C      NZERD10=1
C      NZERD11=1
C      NZERD12=1
C      NZERD13=1
C      NZERD14=1
C      NZERD15=1
C      NZERD16=1
C      NZERD17=1
C      NZERD18=1
C      NZERD19=1
C      NZERD20=1
C      NZERD21=1
C      NZERD22=1
C      NZERD23=1
C      NZERD24=1
C      NZERD25=1
C      NZERD26=1
C
C      GIONEUT 3
C      GIONEUT 4
C      GIONEUT 5
C      GIONEUT 6
C      GIONEUT 7
C      GIONEUT 8
C      GIONEUT 9
C      GIONEUT 10
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C      GIONEUT 53
C      GIONEUT 54
C      GIONEUT 55
C      GIONEUT 56
C      GIONEUT 57
C      GIONEUT 58
C      GIONEUT 59

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```

      EDDYL=EDY
      SCY=SCW(K)
      D0Z=1.0/(DZ*DZ)
      DZ2=DZ*DZ/CC2
      I=1,J=1,L=1,GR=1,LRB, GO TO 15
      TITLE=0.
      WBAR=0.
      DO 10 L=1,KSPCI
      EN=N(J,L)
      WBAR=WBAR+MASS(L)*EN
      TITLE=TITLE+EN
      CONTINUE
      WBAR=WBAR, TITLE
      GAMBA=GAMBA
      GAMBA=1.0-(DNTD*GRAV*GCRN*WBAR)/(BOLTZ*T)*DZ
      DTDZ=DELT*DZDZ
      15 IF(J-EQ.1) GO TO 20
      AJ=DTDZ*(PJ*UTKIN*EUDY)
      BU=1.0-DTDZ*(PU*(DZ+DTKTM*GAM)+PJL*DTKTM*DZDZL_
      *EUDY*GAM)+(PJ*ECD*DZDZL)+(REMV(J)+PJ*SENU*DELT*DZL)*DELT_
      CJ=DTDZ*(PJ*(PJL*DZ+DTKTM*GAM*L*DZDZL)+EDDYL*GAMBA*DZDZL)
      1 +(PJL*PJL*WBAR*DELT/DZL)*DELT
      GO TO 35
      C COMPUTE LOWER BOUNDARY VALUES OF E AND F USING THE LOWER
      C BOUNDARY CONDITIONS.
      C
      20 EUL=0.
      IF((J.EQ.8).OR.(J.EQ.13).OR.(J.EQ.15).OR.(J.EQ.21)) GO TO 30
      IF((J.EQ.22).OR.(J.EQ.23)) GO TD 30
      FUL=(NZERO(1)+DELT*REMV(1))/(1.0+DELT*REMV(1))
      GO TO 70
      30 FUL=N(1,1)
      GO TO 70
      35 IF(J.EQ.K) GO TO 50
      DENOM=1.0/(BJ-CJ*EUL)
      EUL=AJ*DENOM
      CJ=NZERO(CJ)+DELT*FORM(CJ)
      FUL=(DJ+CJ*FUL)*DENOM
      GO TO 70
      C COMPUTE UPPER BOUNDARY VALUES OF E AND F USING THE UPPER
      C BOUNDARY CONDITIONS.
      C
      50 DJ=NZERO(K)+DELT*FORM(K)
      55 VEL(1)=0.
      IF((J.EQ.4).OR.(J.EQ.5).OR.(J.EQ.9)) GO TO 57
      IF((J.EQ.11)) GO TO 57
      IF((J.GE.12).AND.(J.LE.15)) GO TO 57
      IF((J.EQ.19).OR.(J.EQ.19).OR.(J.EQ.20).OR.(J.EQ.22)) GO TO 57
      IF((J.GE.27).AND.(J.LE.29)) GO TO 57
      CALL VELOCITY(J)
      57 CALL VELOCITY(J)
      IF((J.EQ.11)) GO TO 80
      80 IF(VEL(1).NE.0.0) GO TO 85
      85 SSM=VEL(1)*SENU*DEMS
      FLE=(CJ+CJ*FUL)/(BJ-AJ+(GAM+SS)*(DELTAZ(K)*WGT/(BOLTZ*T*TEMP(K))))
```

```

1      SUBROUTINE EFORU(ICF)
2      C IF(ICF .GT. 0), COMPUTE THE E AND F ARRAYS.
3      C IF(ICF .LT. 0), COMPUTE THE SPECIFIC VELOCITIES.
4
5      C DIMENSION H(1)
6
7      C COMMON N(692,30),CONE(424,26),U(692,26),NZERO(692),UZERO(692)
8
9      C COMMON FORM(692),REMV(692),E(692),F(692),Z(692),
10     DEDY(692),TEMP(692),DTEM0Z(692),DELTAZ(692),SUNSET(692)
11
12     C COMMON GRAV,COSD,SIND,TIME,CXI,RADIUS,ISPECI,KSPECI,NREAC,K,
13       K2,KMINI,ITURB,CXIND0N,SCXI,NREAC2
14
15     C COMMON/TIMES/DAYS,DELT,TMAX
16
17     C COMMON/ALTITUDE/ZBOT,ZTOP,TURB,TURBYAR
18
19     C COMMON/BOUNDARY/VEL(26)
20
21     C COMMON/CONSTANT/PI,RADDEG,DEGRAD,RAOSEC,BOLTZ,SIN2I,ATCON
22
23     C COMMON/ATCONS/ATWT(26),MASS(26),THERM(26),SYMBOL(56),POLAR(26)
24
25     C REAL N,NZERO,MASS
26     C EQUIVALENCE (U(1,21),H)
27
28     C I=ICF
29     C IF(I .GT. 0) GO TO 1
30     C I=-I
31
32     1   WGT=MASS(I)
33     GMOVEK=GRAV*WGT/BOLTZ
34     DTOKVM=DELT*BOLTZ/WGT
35     DZ=1.5E+04
36     DO 75 J=1,K
37     ZE=Z(J)
38     T=TEMP(J)
39     DZL=DZ
40     DZ=DELTAZ(J)
41     DLNTDZ=DTEM0Z(J)
42     IF(I .GT. 6) GO TO 6
43     DLNTDZ=2.0*DLNTDZ
44     GCRN=1.0/((1.0+ZE/RADIUS)**2)
45     DTKTM=DTKTM
46     DTKTM=DTKQVM*T
47     GAML=GAM
48     GAM=1.0-(THERM(I)*DLNTDZ+GCRN*GMQVK/T)*DZ
49
50     11   SNUWL=SNUW
51     CALL COLFRE(I,J,SNU,SNUL,OMEG)
52     PJ=1.0/(1.0+SNU*OMEG*DELT)
53     GAM=GAM-H(J)*DZ
54     PJL=PJ
55
56     UJL=UJ
57     UJ=UZ-E(J)
58     IF(ICF .LT. 0) GO TO 60

```

```

1      SUBROUTINE CO(J,RHO,NTOTAL,MBAR,VO)
C
C THIS SUBROUTINE COMPUTES THE TOTAL MASS DENSITY, RHO, THE TOTAL
C CONCENTRATION, NTOTAL, THE MEAN MOLECULAR WEIGHT, MBAR, AND THE
C MEAN MASS VELOCITY, VO, AT ALTITUDE WHERE THE INDEX IS J.
5
C
C COMMON N(692,30),CONC(424,28),U(692,26),NZERO(692),UZERO(692)
C
C COMMON FORM(692),REMY(692),E(692),Z(692),
10    DE(692),TEMP(692),STEMDZ(692),DELTAZ(692),SUNSET(692)
C
C COMMON GRAV,COSD,SIND,TIME,CXI,RADIUS,ISPECI,KSPECI,NREAC,K,
15    K2,KMIN1,ITURB,CKINCON,SCXI,NREAC2
C
C COMMON, VTCO4S/ATWT(26),MASS(26),THERM(26),SYMBOL(56),POLAR(26)
20
C      REAL N,MBAR,NTOTAL,MASS
C
C      RHO=0,
C      NTOTAL=0,
C      RHOC=0.
25
DO 5 M=1,KSPECI
EN=N(J,M)
ENM=MASS(M)*EN
RHC=RHO+ENM
NTOTAL=NTOTAL+EN
IF(M.EQ.21) GO TO 5
RHOC=RHOC+ENM*U(J,M)
CONTINUE
30
MBAR=RHO/NTOTAL
VO=RHOC/RHO
RETURN
END

```

```

1      SUBROUTINE SMOOTH(A,N,M,B)
2
3      C THIS SUBROUTINE SMOOTH'S ARRAY A OF N POINTS WITH AN M POINT
4      C RUNNING MEAN. THE SMOOTHED ARRAY IS STORED IN B.
5
6      C
7      C DIMENSION A(1),B(1)
8      C
9      C IF(M .EQ. 0) GO TO 10
10     C I=(M-1)/2
11     C K=N-M+1
12     C S=M
13     DO 2 J=1,K
14     S=0.
15     C JJJ=J+M-1
16     DO 1 L=J,JJJ
17     S=S+A(L)
18     1 JAY=J+I
19     2 B(JAY)=S/C
20     C I1=0
21     C K=M-2
22     C=C
23     S=0.
24     DO 4 J=1,K
25     S=S+A(J)
26     JAY=I-I1
27     B(JAY)=S/C
28     K=K-2
29     I1=I1+1
30     C IF(K .NE. 1) GO TO 3
31     C I1=1
32     C K=N-M+3
33     C=C-2.
34     C
35     DO 6 J=K,N
36     S=S+A(J)
37     JAY=N-I+I1
38     B(JAY)=S/C
39     K=K+2
40     C=C-2.
41     I1=I1+1
42     C
43     IF(K .NE. N) GO TO 5
44     B(1)=A(1)
45     B(N)=A(N)
46     RETURN
47
48     10 DO 20 J=1,N
49     B(J)=A(J)
50     20 CONTINUE
51     RETURN
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      RETURN
      C   ADJUST THE LARGEST NEGATIVE SPECIES CONCENTRATIONS TO
      C   INSURE CONSERVATION OF CHARGE.
      C
      C
      25    DO 50 J=1,K2
             SUMP=0.
             SUMMA=0.
      65    DO 30 I=1,6
             SUMP=SUMP+N(J,I)
      30    CONTINUE
             DO 35 I=28,29
             SUMP=SUMP+N(J,I)
      35    CONTINUE
             DO 40 I=9,24
             SUMP=SUMP+CONC(J,I)
      40    CONTINUE
             BIG=N(J,30)
      75    DO 70 I=1,8
             IF(CONC(J,I) .LE. BIG) GO TO 70
             BIG=CONC(J,I)
      80    JBIG=+1
      70    CONTINUE
             DO 75 I=1,9
             IF(I .EQ. JBIG) GO TO 75
             IF(I .EQ. 1) SUMMA=SUMMA+N(J,30)
             IF(I .NE. 1) SUMM=SUMM+CONC(J,I-1)
      75    CONTINUE
             BAL=SUMP-SUMM
      90    IF(BAL .LT. 0.) STOP 7
             IF(JBIG .EQ. 1) N(J,30)=BAL
             IF(JBIG .NE. 1) CONC(J,JBIG-1)=BAL
      50    CONTINUE
             KK=K2+1
             DO 65 J=KK,K
             SUMP=0.
      95    DO 55 I=1,6
             SUMP=SUMP+N(J,I)
      55    CONTINUE
             DO 60 I=28,29
             SUMP=SUMP+N(J,I)
      60    CONTINUE
             N(J,30)=SUMP
      65    CONTINUE
             RETURN
      END
      105

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1      SUBROUTINE SMOOTHY (NUM,NSPEC)
C THIS SUBROUTINE SMOOTH SPECIFIED PROFILES WITH AN "IRUN" POINT
C RUNNING MEAN.
5      C
C      DIMENSION NSPEC(1),ZNEW(792),DENS(792),SMUTHN(1)
C
C      COMMON N(692,30),CONC(424,26),U(692,26),NZERO(692),UZERO(692)
C
C      COMMON FORM(692),REMV(692),E(692),F(692),Z(692),
2      DEDY(692),TEMP(692),DTENDZ(692),DELTAZ(692),SUNSET(692)
C
C      COMMON GRAV,COSD,SIND,TIME,CX1,RADIUS,ISPEC1,KSPEC1,NREAC1,K,
1      K2,KMIN1,ITURB,CXINCON,SCX1,NREAC2
C
C      COMMON/ATCONS/ATWT(26),MASS(26),THERM(26),SYMBOL(56),POLAR(26)
C
C      COMMON/CONSTAN/PI,RADdeg,DEGRAD,RADSEC,BOLTZ,SIN2I,ATCON
C
20     C
C      REAL N,NZERO,MASS
C
C      EQUIVALENCE (NZERO,DENS),(FORM,SMUTHN)
C      EQUIVALENCE (E,ZNEW)
C
25     C      DATA IRUN/7/
C
C      IF (NUM .EQ. 0) RETURN
C
C      GOVKT=(GRAV*(1.0/((1.0+400./RADIUS)**2)))/(BOLTZ*TEMP(K))
30     R=6.35107E+08
DO 5   J=1,692
      ZNEW(J)=Z(J)
      CONTINUE
      5   DO 6   J=693,792
      ZNEW(J)=ZNEW(692)+FLDAT(J-692)
      6   CONTINUE
      DO 15  I=1,NUM
      L=NSPEC(I)
      DO 7   J=1,692
      DENS(J)=N(J,L)
      7   CONTINUE
      C
C      EXTRAPOLATE THE CONCENTRATION OF SPECIES L TO 500 KM ASSUMING
C      DIFFUSIVE EQUILIBRIUM ABOVE THE UPPER BOUNDARY.
45     C
C      DO 8   J=693,792
      DENS(J)=DENS(692)*EXP(-(GOVKT*MASS(L)*(ZNEW(J)-400.)*1.0E-5)
1      /((400.*1.0E5/R)*1.0)*((ZNEW(J)*1.0E5/R)+1.0))
      8   CONTINUE
      CALL SMOOTH(DEN,792,IRUN,SMUTHN)
      DO 10  J=1,K
      N(J,L)=SMUTHN(J)
      10    CONTINUE
      15    CONTINUE
      DO 20  I=1,NUM
      L=NSPEC(I)
      LT=7
      20    CONTINUE
      25    CONTINUE
      5

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228  FORMAT(1HO,• DAY IS •,L1,5X,•NIGHT IS •,L1,5X,•DAY-NIGHT•,      GIONEUT
515   1• TRANSITION HEIGHT IS •,F7.2)                                GIONEUT 516
      IF(END) GO TO 270                                              GIONEUT 517
      CALL MOVLEV(U,OLDU,U,17992)                                       SIONEUT 518
      CALL MOVLEV(N,OLDN,N,20760)                                       GIONEUT 519
      CALL MOVLEV(CONC,OLDCON,11024)                                     GIONEUT 520
      CALL SECOND(T1)                                                 GIONEUT 521
      GO TO 1000                                                       GIONEUT 522
      GO TO 1000                                                       GIONEUT 523
      GO TO 1000                                                       GIONEUT 524
      GO TO 1000                                                       GIONEUT 525
      GO TO 1000                                                       GIONEUT 526
      GO TO 1000                                                       GIONEUT 527
      GO TO 1000                                                       GIONEUT 528
      GO TO 1000                                                       GIONEUT 529
      GO TO 1000                                                       GIONEUT 530
      GO TO 1000                                                       GIONEUT 531
      GO TO 1000                                                       GIONEUT 532
      GO TO 1000                                                       GIONEUT 533
      GO TO 1000                                                       GIONEUT 534
      GO TO 1000                                                       GIONEUT 535
      GO TO 1000                                                       GIONEUT 536
      GO TO 1000                                                       GIONEUT 537
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      GO TO 1000                                                       GIONEUT 539
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      GO TO 1000                                                       GIONEUT 542
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      GO TO 1000                                                       GIONEUT 545
      GO TO 1000                                                       GIONEUT 546
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      GO TO 1000                                                       GIONEUT 548
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      GO TO 1000                                                       GIONEUT 552
      GO TO 1000                                                       GIONEUT 553
      GO TO 1000                                                       GIONEUT 554
      GO TO 1000                                                       GIONEUT 555
      GO TO 1000                                                       GIONEUT 556
      GO TO 1000                                                       GIONEUT 557
      GO TO 1000                                                       GIONEUT 558
      GO TO 1000                                                       GIONEUT 559
      GO TO 1000                                                       GIONEUT 560
      GO TO 1000                                                       GIONEUT 561
      GO TO 1000                                                       GIONEUT 562
      GO TO 1000                                                       GIONEUT 563
      GO TO 1000                                                       GIONEUT 564
      GO TO 1000                                                       GIONEUT 565
      GO TO 1000                                                       GIONEUT 566
      GO TO 1000                                                       GIONEUT 567
      GO TO 1000                                                       GIONEUT 568
      GO TO 1000                                                       GIONEUT 569

C ENTER HERE WHEN TIME IS RUNNING OUT OR SENSE SWITCH ONE IS ON.
C
230  CXI=OLDCXI
      SCXI=CXI*DEGRAD
      TIME=TIME-DELT
      DELT=OLDELT
      ZTRANS=OLDZTR
      IF(PRINT) TIMEOUT=TIMEOUT-OPRINT
      PRINT=.TRUE.
      END=.TRUE.
      LM=1
      GO TO 245
      C ENTER HERE ON CALCULATION OF A NEGATIVE CONCENTRATION.
C
245  KFAIL=2
      J=K
      GO TO 255
250  KFAIL=3
      J=L
      WRITE(6,256)KFAIL,SYMBOL(I),Z(J),NEWN,N(J,I),TIME,DELT,SCXI,
      1,TRATE
255  FORMAT(I3,A10,0PF7.1,1P5E20.12,15)
      C ENTER HERE ON ERROR FROM SUBROUTINE CHEMION.
C
254  TIME=OLDTIME
      END=.FALSE.
      CXI=OLDCXI
      DELT=DELT*0.5
      IF(RESET) GO TO 257
      IF(DELT .LT. 2.0*•(-40)) 265,258
      257  RESET=.FALSE.
            TIMEOUT=TIMEOUT-OPRINT
            LM=2
            NIGHT=LASTNT
            DAY=LASTDA
            CALL MOVLEV(OLDN,N,20760)
            CALL MOVLEV(OLDCON,CONC,11024)
            CALL MOVLEV(OLDU,U,17992)
            IF(LM .EQ. 1) 162,1000
            265  WRITE(6,345) TIME
            345  FORMAT(30H THE INCREMENT IS VANISHING AT1PE12.5,5H SEC.)
            GO TO 230
            CONTINUE
            END
      270  CONTINUE
      END

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      WRITE(6,340)
      FORMAT(1H1)
      WRITE(6,196) TIME,SCXI,IFRAME,IDAY,(OUTP(J),J=1,4)
      WRITE(6,310)
      FORMAT(1*, ALT(KM) NO./O2+ RATIO MEAN MOL WT MASS DENSITY*,
     1* TOTAL NUMBER MEAN MASS VEL K-EDDY*)
      DO 325 J=1,K,12
      CALL CO(J,RHO,NTOTAL,MBAR,VMEAN)
      MBAR=MBAR/ATCON
      RNOTO2=N(J,5)/N(J,4)
      WRITE(6,330) Z(J),RNOTO2,MBAR,RHO,NTOTAL,VMEAN,DODY(J)
      CONTINUE
      CALL CO(KMIN1,RHO,NTOTAL,MBAR,VMEAN)
      MBAR=MBAR/ATCON
      RNOTO2=N(KMIN1,5)/N(KMIN1,4)
      WRITE(6,330) Z(KMIN1),RNOTO2,MBAR,RHO,NTOTAL,VMEAN,DODY(K)
      MBAR=MBAR/ATCON
      RNOTO2=N(K,5)/N(K,4)
      WRITE(6,330) Z(K),RNOTO2,MBAR,RHO,NTOTAL,VMEAN,DODY(K)
      FORMAT(10PF8.2,1PE13.5,OPF10.3,4(3X,1PE13.5))
      C
      C PRINT LIST OF CHEMICAL REACTIONS WITH THE RATE OF EACH AT
      C SELECTED ALTITUDES.
      C
      CALL CHEMPR(1)
      CALL CHEMPR(253)
      CALL CHEMPR(337)
      CALL CHEMPR(457)
      CALL CHEMPR(K)
      DELT=AMIN1(2.0*DELT,TMAX)
      IF(.NOT. PRINT) GO TO 161
      C
      C WRITE INTERRUPT FILE ON TAPE 4.
      C
      WRITE(4) TODAY,IFRAME,TIME,SCXI,CXI,DELT,TIMOUT
      DO 2224 L=1,KSPECI
      WRITE(4) (U(U,L),J=1,K)
      CONTINUE
      DO 280 L=1,ISPECI
      IF(L.GT. 30) GO TO 281
      WRITE(4) (N(U,L),J=1,K)
      GO TO 280
      WRITE(4) (CONC(U,L-30),J=1,K2)
      CONTINUE
      280
      ENDFILE 4
      CALL CLOCK(A1)
      WRITE(6,315) A1,IFRAME
      FORMAT(1H0,A17,* FRAME NO. * 14,* HAS BEEN WRITTEN ON THE*,
     1* INTERRUPT TAPE *)
      161
      CALL HOUR(TIME,OUTP)
      CALL CLOCK(A1)
      WRITE(6,227) A1,TIME,DELT,SCXI,ITRATE,(OUTP(J),J=1,4),DTIME
      FORMAT(1H0,A10,* TIME **, PE12.5 * SEC. TIME INCREMENT **,
     1E12.5,* CH1 **, OPF7.2,* ITERATIONS **, 14,* HOUR **,4A1,
     2* CP TIME **, F10.2,* SEC *)
      WRITE(6,228) DAY,NIGHT,ZTRANS

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400      1*NUMBER CHEMISTRY CHEMISTRY*, ALT NUMBER *.
        2*CHEMISTRY (CHEMISTRY*)
201      FFORMAT(*,KM,DENSITY FORMATION REMOVAL*,BX,*.
        1*DENSITY FORMATION REMOVAL*,BX,*.
        KM DENSITY*.
2* FORMATION REMOVAL*)
        WRITE(6,200)
        WRITE(6,201)
        WRITE(6,202) (Z(J),CONC(J,1-3),E(J),H(J),Z(J),CONC(J,1-2),F(J),
1          UZERO(J),Z(J),CONC(J,1-1),FORM(J),REMV(J),J=1,K2,12)
        WRITE(6,202) Z(K2),CONC(K2,1-3),E(K2),H(K2),Z(K2),CONC(K2,1-2),
1          F(K2),UZERO(K2),Z(K2),CONC(K2,1-1),FORM(K2),REMV(K2)
202      FORMAT(3(0PF7.2,1P3E10.3,6X))
210      CONTINUE
I=25
        DO 237 LL=1,2
        DO 211 J=1,K2
        CALL CHEM(I+30,J)
        REMV(J)=REMV(J)*CONC(J,I)
211      CONTINUE
        IF(LL.EQ. 2) GO TO 240
        DO 236 J=1,K2
        NZERO(J)=FORM(J)
        UZERO(J)=REMV(J)
236      CONTINUE
        I=I+1
237      CONTINUE
240      DO 241 J=1,K2
        E(J)=0.
        F(J)=0.
        H(J)=0.
        DG 312 I=1,6
        E(J)=E(J)+N(J,I)
        E(J)=E(J)+N(J,28)+N(J,29)
        DO 313 I=9,24
        E(J)=E(J)+CONC(J,I)
238      CONTINUE
        DO 314 I=1,8
        F(J)=F(J)+CONC(J,I)
239      CONTINUE
        H(J)=F(J)/N(J,30)
240      CONTINUE
        DG 314 I=1,8
        F(J)=F(J)+CONC(J,I)
241      CONTINUE
        WRITE(6,340)
        TIME,SCX1,IFRAME,1DAY,(OUTP(J),J=1,4)
        WRITE(6,196) SYMBOL(55),SYMBOL(56)
212      FORMAT(10X,A10,* PROFILE*,20X,A10,* PROFILE*)
        WRITE(6,213)
213      FORMAT(* ALT NUMBER CHEMISTRY CHEMISTRY ALT NUMBER *
        1*CHEMISTRY CHEMISTRY TOTAL IONS TOTAL IONS LAMBDAA*)
        WRITE(6,214)
214      FORMAT(* KM DENSITY FORMATION REMOVAL KM DENSITY *
        1* FORMATION REMOVAL POSITIVE NEGATIVE*)
        WRITE(6,215) (Z(J),CONC(J,25),NZERO(J),Z(J),CONC(J,26),
1 FORM(J),REMV(J),E(J),F(J),H(J),J=1,K2,12)
        WRITE(6,215) Z(K2),CONC(K2,25),NZERO(K2),UZERO(K2),Z(K2),
        CONC(K2,26),FORM(K2),REMV(K2),E(K2),F(K2),H(K2)
215      FORMAT(0PF7.2,1P3E10.3,0PF7.2,1P3E10.3,3E12.5)

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C COMPUTE EDDY DIFFUSION VELOCITIES.
C
C     BIG=N(J+1,I)/EN
C     IF(EN .GT. 1.0E-10) GO TO 165
C     IF(BIG .GT. 100.) GO TO 170
C     165    DED(J)=-(DEOY(J)/DETAZ(J))•(BIG-1.0+GAMA•DETAZ(J))
C     170    CONTINUE
C
C     COMPUTE FLUX GRADIENTS.
C
C     DEDZ(1)=0.
C     DFDZ(1)=0.
C     DO 175 J=2,K
C     DEDZ(J)=0.
C     DFDZ(J)=N(J,I)*U(J,I)-N(J-1,I)*U(J-1,I)/DETAZ(J)
C     IF(J .GT. 1)URB) GO TO 175
C     DEDZ(J)=N(J,I)*DED(J)-N(J-1,I)*DED(J-1)/DETAZ(J)
C     175    CONTINUE
C
C     COMPUTE CHEMISTRY PROFILES.
C
C     DO 180 J=1,K
C     CALL CHEM(I,J)
C     REMV(J)=REMV(J)*N(J,I)
C     180    CONTINUE
C
C     COMPUTE THE TOTAL FLUX FOR EACH SPECIES.
C
C     DO 181 J=1,K
C     PHI(J)=(U(J,I)+DED(J))*N(J,I)
C     181    CONTINUE
C     CALL HOUR(TIME,OUTP)
C     WRITE(6,335) SYMBOL(I)
C     FORMAT(1H1,60X,A10,* PROFILE*)
C     335    WRITE(5,390) TIME,DELT,SCX,ITER(I),IFRAME,IDAY,(OUTP(J),J=1,4)
C
C     390    FORMAT(1H,* TIME =*,1PE12.5,* SEC. TIME INCREMENT =*,E12.5,
C           1* CHI **,OPF7.2,* DEGREES ITERATIONS **,13,* FRAME**,
C           214,* DAY **,13,* HOUR **,4A1)
C     WRITE(6,295)
C     295    FORMAT(* ALT      NUMBER      SPECIFIC      SPECIFIC*,
C           1* TURBULENT   TURBULENT   CHEMISTRY      CHEMISTRY*,
C           2* TOTAL      DIFFUSION*)      REMOVAL
C
C     300    FORMAT(* KM      DENSITY      VELOCITY      GRADIENT      VELOCITY*,
C           1,* GRADIENT FORMATION      FLUX*,
C           2,* VELOCITY*)      REMOVAL
C
C     330    WRITE(6,300)
C
C     330    FORMAT(* KM      DENSITY      VELOCITY      GRADIENT      VELOCITY*,
C           1,* GRADIENT FORMATION      FLUX*,
C           2,* VELOCITY*)      REMOVAL
C
C     330    WRITE(6,320) (Z(J,I),U(J,I),DFDZ(J),DED(J),FORM(J),FORM(J),
C           1PEMV(J),PHI(J),VBAR(J),J=1,K,12)
C     330    FORMAT(0PF8.2,1PF8.13.5)
C     330    WRITE(6,320) Z(K),N(K,I),U(K,I),DFDZ(K),DED(K),FORM(K),
C           1REMV(K),PHI(K),VBAR(K)
C
C     195    CONTINUE
C     L=1
C     341    I=27,30
C     342    I=191,J=1,K
C     343    CALL CEM(I,J)
C
C     287    GONEUT
C     288    GONEUT
C     289    GONEUT
C     290    GONEUT
C     291    GONEUT
C     292    GONEUT
C     293    GONEUT
C     294    GONEUT
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C     298    GONEUT
C     299    GONEUT
C     300    GONEUT
C     301    GONEUT
C     302    GONEUT
C     303    GONEUT
C     304    GONEUT
C     305    GONEUT
C     306    GONEUT
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C     320    GONEUT
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C     323    GONEUT
C     324    GONEUT
C     325    GONEUT
C     326    GONEUT
C     327    GONEUT
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C     338    GONEUT
C     339    GONEUT
C     340    GONEUT
C     341    GONEUT
C     342    GONEUT
C     343    GONEUT

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      IF(LITMAX .EQ. 1) GO TO 139
      IF(LITRATE .EQ. 1) GO TO 130
      C DETERMINE WHICH SPECIES REQUIRE FURTHER ITERATION.
      C
      BIG=0.
      DO 120 I=1,KSPECI
      IF(INWHICH(I) .EQ. 0) GOTO 120
      IF(BIGEST(I) .GT. ERSPECI) GO TO 117
      IWHICH(I)=0
      ITER(I)=ITERATE
      IF(BIGEST(I) .LT. BIG) GO TO 120
      BIG=BIGES(I)
      117
      120 CONTINUE
      C
      ERS=ERSMAX
      IF(BIG .LE. ERS) GO TO 139
      130 CONTINUE
      139 DO 135 I=1,KSPECI
      IF(INWHICH(I) .EQ. 0) GO TO 135
      ITER(I)=ITERATE
      135 CONTINUE
      CALL SECOND(12)
      DTIME=12-T1
      C
      C IF THIS IS A PRINT STEP, COMPUTE FLUXES AND FLUX GRADIENTS FOR OUTPUT
      C
      IF(PRINT) GO TO 162
      IF(TIME .LT. (DAYS*8.64E+04)) GO TO 160
      END=.TRUE.
      160 DAY=(ITIME*4.32E+04)/8.65E+04)+1
      IFRAME=IFRAME+1
      DO 163 J=1,K
      CALL RATECN(J)
      CALL MOVLEV(DK(128),RATES(1,J),88)
      163 CONTINUE
      C
      C COMPUTE MOLECULAR AND EDDY DIFFUSION FLUXES.
      C
      DO 195 I=1,KSPECI
      IF(I .EQ. 21) GO TO 195
      DO 170 J=1,K
      C
      C COMPUTE MOLECULAR DIFFUSION VELOCITIES.
      C
      CALL CD(J,RH0,NTOTAL,MBAR,VMEAN)
      VBAR(J)=0.
      EN=N(.1.)
      DED(J)=0.
      IF(EN .EQ. 0.) GO TO 170
      VBAR(-1)=EV(J,I)-VMEAN
      IF(J .LT. ITURB) GO TO 170
      T=TEND(J)
      ZZ=Z(J)
      GAMA=2*EV(2*(J)+1,GRA)*GRA*CCP(22)*MBAR/(BOLTZ*T)
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      CALL EFORU(I)
C COMPUTE HEIGHT PROFILES AT (T+DELTA T) USING THE RECURSION
C RELATION N(J)=N(J+1)*E(J)+F(J).
C
C NEWN=F(K)
C
C   IF(NEWN .GT. 0.) GO TO 10
C   IF(N(K,1) .GT. 1.0E-02) GO TO 245
C   NEWN=N(K,1)
C   10   IF(ITERATE .EQ. 1) GO TO 70
C   IF(NEWN .LT. 1.0) GO TO 70
C   RELER=ABS((NEWN-N(K,1))/NEWN
C   BIGEST(I)=RELER
C   N(K,1)=NEWN
C   DO 100 J=1,KMINI
C     L=K-J
C     NEWN=E(L)*N(L+1,1)+F(L)
C     IF(NEWN .GT. 0.) GO TO 11
C     IF(N(L,1) .GT. 1.0E-02) GO TO 250
C     NEWN=N(L,1)
C     11   IF(ITERATE .EQ. 1) GO TO 95
C     IF(NEWN .LT. 1.0) GO TO 95
C     RELER=ABS((NEWN-N(L,1))/NEWN
C     IF(BIGEST(I) .LT. RELER) BIGEST(I)=RELER
C     95   N(L,1)=NEWN
C     100 CONTINUE
C     105 CONTINUE
C
C   COMPUTE THOSE SPECIES TO BE DETERMINED FROM CHEMISTRY ONLY.
C   CALL CHEMION
C
C   IF(END) GO TO 254
C
C   GO GET THE VELOCITIES FOR SPECIES IN THE TRANSPORT MODE.
C   DO 110 I=1,KSPECI
C   IF(I .EQ. 21) GO TO 110
C   CALL MOVLEV(OLDDU(1,1),UZERO,K)
C   CALL CHEM(I,KMINI)
C   CALL CHEM(I,K)
C   IF(I .NE. 1) GO TO 108
C   DO 107 J=2,K
C     H(J)=(1.0-(N(J-1,30)/N(J-30)))/DELTAZ(J-1)
C   107 CONTINUE
C     H(1)=H(2)
C   108 CALL EFORU(-1)
C
C   TEST FOR TERMINATION OF CALCULATIONS.
C   CALL SSWITCH(NSWITCH,NSTATUS)
C   IF(NSTATUS.NE.2) GOTO 230
C   CALL SECOND(TA)
C   IF((TTYPE-TA).LT.TOFF) GOTO 230
C   CONTINUE
C
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      GO TO 625 J=1,50
      WRITE(6,610),DIS(15,1),DIS(16,1),DIS(17,14)
      CONTINUE
      610
      CONTINUE FA SE:
24      IF(LTTRATE .NE. 1) GO TO 6
      IF(LTTRATE .NE. 11) GO TO 6
      IF(.TRUE.) EQ. 01 GO TO 6
      IF(.TRUE.) EQ. 01 GO TO 6
      CALL SMOOTH(NUM,NSPEC)
      WRITE(6,2)
      2      FORMAT(1HO,* THE FOLLOWING SPECIES PROFILES HAVE BEEN*)
      1      * SMOOTHED.*)
      DO 4 I1=1,NUM
      L=NSPEC(I1)
      WRITE(6,3) SYMBOL(L)
      3      FORMAT(10X A10)
      FORMAT(10X A10)
      CONTINUE
      4      RECOMPUTE THE VELOCITY PROFILES FOR THOSE SPECIES WHOSE
      CONCENTRATION PROFILES HAVE BEEN SMOOTHED.
      C
      35      IP=0
      DO 29 IM*1,NUM
      29      I=NSPEC(IM)
      40      IP=IP+1
      41      IF(IP .NE. 11) GO TO 40
      CALL MOVLEY(OLDU(1,11),UZERO,K)
      CALL CHEM(11,KMINI)
      CALL CHEM(11,K)
      IF(IM .NE. 1) GO TO 28
      DO 27 J=2,K
      H(J)=(1.0-(N(J-1,30),N(J,30)))/DELTAZ(J-1)
      CONTINUE
      27      H(1)=H(2)
      28      CALL EFORU(-11)
      29      CONTINUE
      150      CALL MOVLEY(U,OLDU,17992)
      CALL MOVLEY(N,OLDN,20760)
      CALL MOVLEY(CONC,OLDCON,11024)
      6      1F(1WHICH(1) .EQ. 0) GO TO 105
      CALL MOVLEY(OLDN(1,1),NZERO,K)
      CALL MOVLEY(OLDU(1,1),UZERC,K)
      155      C GO GET CHEMISTRY FORMATION AND REMOVAL RATES FOR SPECIES 1.
      C
      160      DO 30 J=1,K
      30      CALL CHEM(1,J)
      C TEST FOR TERMINATION OF CALCULATIONS.
      C
      165      CALL SSWATCH(NSWITCH,INSTATUS)
      1F(INSTATUS .NE. 2) GO TO 230
      CALL SECOND(TUSED)
      1F((TIME-TUSED) .LT. TOFF) GO TO 230
      30      CONTINUE
      C GO SET E AND F ARRAYS.
      C
      170
      171
      172
      GIONEUT 116
      GIONEUT 117
      GIONEUT 118
      GIONEUT 119
      GIONEUT 120
      GIONEUT 121
      GIONEUT 122
      GIONEUT 123
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      GIONEUT 171
      GIONEUT 172

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      GIONEUT 5
      GIONEUT 6
      GIONEUT 61
      GIONEUT 62
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      GIONEUT 110
      GIONEUT 111
      GIONEUT 112
      GIONEUT 113
      GIONEUT 114
      GIONEUT 115

      C CALL X1.CAL
      C CALL SSWTCH,NSTATUS,144
      C LASTR=21445
      C ASNT=NIGHT
      LASTDA=DAY
      CALL TIMER
      QDDELT=DELT
      IF(.NOT. DAY .OR. NIGHT) JLOOK=.TRUE.
      IF(NIGHT) JLOOK=.FALSE.
      SCX1=CX1*DEGRAD
      DO 20 I=1,KSPECI
      INHICH(I)=1
      ITER(I)=999
      20 CONTINUE
      INHICH(21)=0

      C BEGINNING OF ITERATION LOOP.
      C
      DO 130 ITRATE=1,ITMAX
      DO 105 I=1,KSPECI
      BIGEST(I)=0.
      IF(I .NE. 1) GO TO 24
      DO 22 J=2,K
      H(J)*(1.0-(N(J-1,30)/N(J,30)))/DELTAZ(J-1)
      22 CONTINUE
      H(1)=H(2)

      C TEST FOR TERMINATION OF CALCULATIONS.
      C
      CALL SSWTCH(NSTATUS,NSTATUS)
      IF(INSTATUS .NE. 2) GO TO 230
      CALL SECOND(TUSED)
      IF((TYME-TUSED) .LT. TOFF) GO TO 230
      DO 21 J=1,K
      CALL RATECN(J)
      CALL MOVLEV(DK(12B),RATES(1,J),88)
      21 CONTINUE
      100
      600 FORMAT(1H1,10X,* CHI **,F7.2,* DEGREES.*)
      105
      WRITE(6,600) SCXI
      605 FORMAT(1H0,* ALT (KM) 02 DENSITY 02 COLUMN 03 DENSITY*.
      1* 03 COLUMN 0J02 K192 0J02 K193 0J03 K194 *.
      2*0J03 K195 0JH20 K196 0JH202 K197*)
      DD 615 J=1,59
      WRITE(6,610) ((DIS(LL,J),LL=1,19),(DIS(LL,J),LL=1,6))
      610 FORMAT(1P10.2,X,1P10E12.5)
      615 CONTINUE
      WRITE(6,600) SCXI
      WRITE(6,620) SCXI
      620 FORMAT(1H0,* ALT (KM) 0C02 K198 0J02 K199 0J02 K200*.
      * 0J03 K201 0J02 K202 0J03 K203 0J02 K204 0J03 K205*)

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115      1      -CJ•EUL)
          EUL=0.
          GO TO 70
          FUL=(NZERO(X)+FORM(X)•DELT)/(1.0+RENV(X)•DELT)
120      EUL=0.
          GO TO 70
          C   COMPUTE THE SPECIFIC VELOCITY ARRAY FOR SPECIES ICF.
          C
          IF(J .EQ. K) GO TO 56
          AJ=N(J+1,1)/N(J,1)
          IF(N(J,1) .GT. 1.0E-10) GO TO 61
          IF(AJ .LT. 100.) GO TO 61
          U(J,1)=0.
          GO TO 75
          U(J,1)=PJ*(UJ-(DTKTM/DZ)*(AJ-GAM)+DELT*SNW)
130      GO TO 75
          U(K,1)=VEL(1)
          GO TO 75
          E(J)=EUL
          F(J)=FUL
          CONTINUE
          RETURN
          END

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      SUBROUTINE COLFRE(I,J,SNU,SNUW,OMEG)
C
C THIS SUBROUTINE COMPUTES THE SUM OF THE COLLISION FREQUENCIES
C AND THE SUM OF THE COLLISION FREQUENCIES TIMES THE VELOCITIES.
C SEE BANKS AND KOCHARIS, "AERONOMY", CHAPTER 9.
C
C COMMON N(692,30), CONC(424,26), U(692,26), NZERO(692)
C
C COMMON FORM(692), REV(692), E(692), F(692), Z(692),
C DEDY(692), TEMP(692), DTENDZ(692), DELTAZ(692), SUNSET(692)
C
C COMMON GRAY, COSD, SIND, TIME, CX1, RADIUS, ISPECI, KSPECI, NRREC,
C K2, KMIN1, ITURB, CXINCON, SCX1, NPEAC2
C
C COMMON/ATCNS/ATWT(26), MASS(26), THERM(26), SYMBOL(56), POLAR(26)
C
C COMMON/CONSTAN/PI, RADDEG, DEGRAD, RADSEC, BOLTZ, SIN21, ATCON
C
C COMMON/GCOLFRE/D(26,26)
C
C REAL N, MASS
C
C DATA GAUS/Q.25/, CHARG/4.803E-10/
C
C SNU = C,
C SNUW = C,
C OMEG = I, 0
C AMUI = ATWT(I)
C
C T = TEMP(J)
C ST = SQRT(T)
C T1 = 1.0/T * 1.5
C 12 = 300.0/T
C 13 = SQRT(2.0*T)
C DO 50 L=1,KSPECI
C AMUL = ATWT(L)
C PNAS = 1.0/(SORT((AMUI*AMUL)/(AMUI+AMUL)))
C
C IF(L .LE. 6) GO TO 15
C IF(L .GT. 6) GO TO 25
C
C COMPUTE NEUTRAL-NEUTRAL COLLISION FREQUENCIES.
C
C ANU=N(J,L)*D(T,L)*ST
C SNU=SNU+ANU
C IF(L .EQ. 21) GO TO 50
C SNUW=SNUW+ANU*U(J,L)
C GO TO 50
C
C 15 IF(L .GT. 6) GO TO 24
C
C COMPUTE ION-ION COLLISION FREQUENCIES.
C
C ANU=1.3*N(J,L)*RMAS*T1
C GO TO 26
C
C COMPUTE ION-NEUTRAL COLLISION FREQUENCIES.
C
C 25 IF(J .EQ. 17)
C GO TO 26
C
C 26
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24   II=L
27   ALF=POLAR(II)  GO TO 50
60
C   AVERAGE ION-NEUTRAL COLLISION FREQUENCY FOR MOMENTUM TRANSFER.
C   BANKS AND KOCHARTS, EQUATION 9.73.
C
65   ANU=2.6E-09*N(J,L)*SQRT(ALF)*RMAS
      IF(((I.EQ.1).AND.(L.EQ.7)).OR.((I.EQ.7).AND.(L.EQ.1))) GO TO 1
      IF(((I.EQ.1).AND.(L.EQ.11)).OR.((I.EQ.11).AND.(L.EQ.1))) GO TO 11
      IF(((I.EQ.2).AND.(L.EQ.8)).OR.((I.EQ.8).AND.(L.EQ.2))) GO TO 2
      IF(((I.EQ.2).AND.(L.EQ.21)).OR.((I.EQ.21).AND.(L.EQ.2))) GO TO 21
      IF(((I.EQ.2).AND.(L.EQ.23)).OR.((I.EQ.23).AND.(L.EQ.2))) GO TO 23
      IF(((I.EQ.3).AND.(L.EQ.7)).OR.((I.EQ.7).AND.(L.EQ.3))) GO TO 3
      IF(((I.EQ.3).AND.(L.EQ.8)).OR.((I.EQ.8).AND.(L.EQ.3))) GO TO 4
      IF(((I.EQ.3).AND.(L.EQ.11)).OR.((I.EQ.11).AND.(L.EQ.3))) GO TO 5
      IF(((I.EQ.4).AND.(L.EQ.18)).OR.((I.EQ.18).AND.(L.EQ.4))) GO TO 6
      IF(((I.EQ.4).AND.(L.EQ.8)).OR.((I.EQ.8).AND.(L.EQ.4))) GO TO 61
      IF(((I.EQ.6).AND.(L.EQ.8)).OR.((I.EQ.8).AND.(L.EQ.6))) GO TO 7
      IF(((I.EQ.6).AND.(L.EQ.16)).OR.((I.EQ.16).AND.(L.EQ.6)))
         BNU=ANU*(AMUL/(AMUL+ANUL))
         SNU=SNU+BNU
         IF(I.EQ.-L) GO TO 50
         IF(L.EQ.-21) GO TO 50
         SNU=SNU+BNU+U(J,L)
         GO TO 50
26
C   AVERAGE COLLISION FREQUENCY FOR CHARGE EXCHANGE REACTION H+ - O.
1   ENG=3.8E-10*N(J,L)
GO TO 8
85
C   AVERAGE REDUCED CHARGE EXCHANGE COLLISION FREQUENCY FOR H+ - H.
1   ANU=ANU+1.0E-10*T3*N(J,L)
GO TO 26
90
C   AVERAGE COLLISION FREQUENCY FOR CHARGE EXCHANGE REACTION HE+ + O2.
1   ENG=2.0E-10*N(J,L)
GO TO 8
95
C   AVERAGE REDUCED CHARGE EXCHANGE COLLISION FREQUENCY FOR HE+ + HE.
2   ANU=ANU+3.0E-11*T3*N(J,L)
GO TO 26
100
C   AVERAGE COLLISION FREQUENCY FOR CHARGE EXCHANGE REACTION HE+ + N2.
2   ENG=3.5E-10*N(J,L)
GO TO 8
105
C   AVERAGE REDUCED CHARGE EXCHANGE COLLISION FREQUENCY FOR O+ + O.
3   ENG=3.0E-11*T3*N(J,L)
GO TO 26
110
C   AVERAGE CC-COLLISION FREQUENCY FOR CHARGE EXCHANGE REACTION O+ + O2.
31  ANU=ANU+1.6E-11*T3*N(J,L)
GO TO 26
C

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115      C GIONEUT 1009
        C GIONEUT 1010
        C GIONEUT 1011
        C GIONEUT 1012
        C GIONEUT 1013
        C GIONEUT 1014
        C GIONEUT 1015
        C GIONEUT 1016
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        C GIONEUT 1019
        C GIONEUT 1020
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        C GIONEUT 1022
        C GIONEUT 1023
        C GIONEUT 1024
        C GIONEUT 1025
        C GIONEUT 1026
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        C GIONEUT 1036
        C GIONEUT 1037
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        C GIONEUT 1039
        C GIONEUT 1040
        C GIONEUT 1041
        C GIONEUT 1042
        C GIONEUT 1043
        C GIONEUT 1044

    4   ENG=2.0E-11*(T2**0.4)*N(J,L)
    5   GO TO 8
    C AVERAGE COLLISION FREQUENCY FOR CHARGE EXCHANGE REACTION O+ + H.
    C ENG=6.8E-10*N(J,L)
    5   GO TO 8
    C AVERAGE COLLISION FREQUENCY FOR CHARGE EXCHANGE REACTION O2+ + NO.
    C ENG=4.4E-10*N(J,L)
    6   GO TO 8
    C AVERAGE REDUCED CHARGE EXCHANGE COLLISION FREQUENCY FOR O2+ + O2.
    C ANU=ANU+1.1E-11*T3*N(J,L)
    61  GO TO 26
    C AVERAGE REDUCED CHARGE EXCHANGE COLLISION FREQUENCY FOR N+ + N.
    C ANU=ANU+1.7E-11*T3*N(J,L)
    71  GO TO 25
    C AVERAGE COLLISION FREQUENCY FOR CHARGE EXCHANGE REACTION N+ + O2.
    C ENG=5.5E-10*(T2**0.17)*2.0*N(J,L)
    7   ANU=ANU+1.73725*ENG
    8   GO TO 26
    50  CONTINUE
    145 IF(I .GT. 6) RETURN
        G1=((GAUS*CHARG)/(MASS(1)*3.0E10))**2
        G2=SNU**2
        OMEG=(G1+G2)/(G1*SIN2I+G2)
        RETURN
    END
    150

```

```

1          SUBROUTINE VELOCY(I)
2
3          C THIS SUBROUTINE COMPUTES THE VELOCITIES AT THE UPPER BOUNDARY
4          C BY SOLVING OVER THE TOPMOST HEIGHT STEP THE ORDINARY DIFFERENTIAL
5          C EQUATION RESULTING FROM SETTING DN/DT=0 IN THE CONTINUITY EQUATION.
6
7          COMMON N(692, 30), CONC(424, 26), U(692, 26), NZERO(692), UZERO(692)
8
9          COMMON FORM(692), RENV(692), E(692), F(692), Z(692),
10             2 DEDY(692), TEMP(692), DTENDZ(692), DELTAZ(692), SUNSET(692)
11
12          COMMON GRAV, COSD, SIND, TIME, CX1, RADIUS, ISPECI, KSPECI, NRREAC, K,
13             1 K2, KMIN1, ITURB, CXINON, SCXI, NRREAC2
14
15          COMMON/BOUNDRY/VEL(26)
16
17          COMMON/RHS10/Z, AN, BN, AP, BP, AL, BL
18
19          REAL N
20          AF(D1, D2, DZ) = (D1-D2)/DZ
21          BF(D1, D2, D3, D4, DZ) = (D2*D3-D1*D4)/DZ
22
23          ZE=Z(KMIN1)
24          START=U(KMIN1, I)
25          DZ=DELTAZ(K)
26          DI=N(K, I)
27          D2=N(KMIN1, I)
28          D3=Z(K)*1.0E+05
29          D4=Z(KMIN1)*1.0E+05
30          AN=AF(D1, D2, DZ)
31          BN=BF(D1, D2, D3, D4, DZ)
32          DI=FORM(K)
33          D2=FORM(KMIN1)
34          AP=AF(D1, D2, DZ)
35          BP=BF(D1, D2, D3, D4, DZ)
36          D1=REMV(K)*N(K, I)
37          D2=REMV(KMIN1)*N(KMIN1, I)
38          AL=AF(D1, D2, DZ)
39          BL=BF(D1, D2, D3, D4, DZ)
40          CALL ODE(I, START, ZE, DZ)
41          VEL(I)=START
42          RETURN
43          END

```

```

1      SUBROUTINE ODE(N,START,SZ,DZ)
2
3      C   SUBPROGRAM TO SOLVE A SET OF ORDINARY DIFFERENTIAL EQUATIONS OF
4      C   SIZE N USING THE KUTTA-MERSON SCHEME.
5
6      C   DIMENSION V(7,2),START(2),SLOPE(2)
7
8      DATA IPASS/0/
9
10     IF(IPASS.NE.0) GO TO 1
11
12     C1=1.0/3.0
13     C2=1.0/6.0
14     C3=1.0/8.0
15     C4=3.0/8.0
16     C5=1.0/2.0
17     C6=3.0/2.0
18     C7=2.0/3.0
19
20     IPASS=1
21     DO 5 J=1,N
22     V(3,J)=V(1,J)=START(J)
23     CONTINUE
24     ZE=SZ
25
26     DO 55 I1=1,S
27     CALL RHSIZE,SLOPE,N,V
28     DO 6 J=1,N
29     V(2,J)=SLOPE(J)
30     CONTINUE
31     GO TO 10,15,20,25,30,11
32
33     DO 11 J=1,N
34     V(4,J)=V(2,J)*DZ
35     V(1,J)=V(3,J)+C1*V(4,J)
36     CONTINUE
37     GO TO 40
38
39     DO 16 J=1,N
40     V(1,J)=V(3,J)+C2*V(4,J)+C2*V(2,J)*DZ
41     CONTINUE
42     GO TO 55
43
44     DO 21 J=1,N
45     V(5,J)=V(2,J)*DZ
46     V(1,J)=V(3,J)+C3*V(4,J)+C4*V(5,J)
47     V(7,J)=V(1,J)
48     CONTINUE
49     GO TO 50
50
51     DO 31 J=1,N
52     V(1,J)=V(3,J)+C2*V(4,J)+C7*V(5,J)-C6*V(6,J)+2.0*V(6,J)*DZ
53     CONTINUE
54     GO TO 55
55
56     ZE=SZ+C1*DZ+1.0E-05
57
58     DO 30 J=1,N
59     V(1,J)=V(3,J)+C2*V(4,J)+C7*V(5,J)+C2*V(2,J)*DZ
60     CONTINUE
61     GO TO 55
62
63     ZE=SZ+C5*DZ+1.0E-05
64
65     DO 30 J=1,N
66     V(1,J)=V(3,J)+C2*V(4,J)+C7*V(5,J)+C2*V(2,J)*DZ
67     CONTINUE
68     GO TO 55
69
70     ZE=SZ+C1*DZ+1.0E-05
71
72     DO 31 J=1,N
73     V(1,J)=V(3,J)+C2*V(4,J)+C7*V(5,J)+C2*V(2,J)*DZ
74     CONTINUE
75     GO TO 55
76
77     ZE=SZ+C5*DZ+1.0E-05
78
79     DO 30 J=1,N
80     V(1,J)=V(3,J)+C2*V(4,J)+C7*V(5,J)+C2*V(2,J)*DZ
81     CONTINUE
82     GO TO 55
83
84     ZE=SZ+C1*DZ+1.0E-05
85
86     DO 31 J=1,N
87     V(1,J)=V(3,J)+C2*V(4,J)+C7*V(5,J)+C2*V(2,J)*DZ
88     CONTINUE
89     GO TO 55
90
91     ZE=SZ+C5*DZ+1.0E-05
92
93     DO 30 J=1,N
94     V(1,J)=V(3,J)+C2*V(4,J)+C7*V(5,J)+C2*V(2,J)*DZ
95     CONTINUE
96     GO TO 55
97
98     ZE=SZ+C1*DZ+1.0E-05
99
100    DO 31 J=1,N
101    V(1,J)=V(3,J)+C2*V(4,J)+C7*V(5,J)+C2*V(2,J)*DZ
102    CONTINUE
103
104    ZE=SZ+C5*DZ+1.0E-05
105
106    DO 30 J=1,N
107    V(1,J)=V(3,J)+C2*V(4,J)+C7*V(5,J)+C2*V(2,J)*DZ
108    CONTINUE
109
110    ZE=SZ+C1*DZ+1.0E-05
111
112    DO 31 J=1,N
113    V(1,J)=V(3,J)+C2*V(4,J)+C7*V(5,J)+C2*V(2,J)*DZ
114    CONTINUE
115
116    ZE=SZ+C5*DZ+1.0E-05
117
118    DO 30 J=1,N
119    V(1,J)=V(3,J)+C2*V(4,J)+C7*V(5,J)+C2*V(2,J)*DZ
120    CONTINUE
121
122    ZE=SZ+C1*DZ+1.0E-05
123
124    DO 31 J=1,N
125    V(1,J)=V(3,J)+C2*V(4,J)+C7*V(5,J)+C2*V(2,J)*DZ
126    CONTINUE
127
128    ZE=SZ+C5*DZ+1.0E-05
129
130    DO 30 J=1,N
131    V(1,J)=V(3,J)+C2*V(4,J)+C7*V(5,J)+C2*V(2,J)*DZ
132    CONTINUE
133
134    ZE=SZ+C1*DZ+1.0E-05
135
136    DO 31 J=1,N
137    V(1,J)=V(3,J)+C2*V(4,J)+C7*V(5,J)+C2*V(2,J)*DZ
138    CONTINUE
139
140    ZE=SZ+C5*DZ+1.0E-05
141
142    DO 30 J=1,N
143    V(1,J)=V(3,J)+C2*V(4,J)+C7*V(5,J)+C2*V(2,J)*DZ
144    CONTINUE
145

```

GIONEUT 1146
GIONEUT 1147
GIONEUT 1148
GIONEUT 1149
GIONEUT 1150
GIONEUT 1151

55 CONTINUE
 DO 60 J=1,N
 START(J)=V(1,J)
 CONTINUE
 RETURN
 END

60

```

1      SUBROUTINE RHS(ZE,SLOPE,N,V)
2
3      C THIS SUBROUTINE COMPUTES THE RIGHT HAND SIDE OF THE ORDINARY
4      C DIFFERENTIAL EQUATION FOR THE VELOCITY AT THE UPPER BOUNDARY.
5
6      C DIMENSION V(7,2),SLOPE(2)
7
8      C COMMON/RHSIDE/AN,BN,AP,BP,AL,BL
9
10     C FF(D1,D2,D3)=D1*D2*1.0E+05*D3
11
12     C EN=EFF(AN,ZE,BN)
13     C EP=EFF(AP,ZE,BP)
14     C EL=EFF(AL,ZE,BL)
15     C SLOPE(1)=(EP-EL-V(1,1)*AN)/EN
16     C RETURN
17     C END
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1      SUBROUTINE TIMER
2      C THIS SUBROUTINE CONTROLS THE TIME AND THE TIME STEP.
3      C
4      COMMON N(692,30),CONC(424,26),U(692,26),NZERO(69-,UZERO(692)
5      C
6      COMMON FORM(692),REMV(692),E(692),F(692),Z(692),
7      DEDY(692),TEMP(692),DTENMZ(692),DELTAZ(692),SUNSET(692)
8      K2,KMIN1,ITURB,CAXNOON,SXI,NREAC2
9      COMMON/TIMES/DAYS,DELT,TMAX,TSET,TRISE
10     COMMON/OJPUTS/TIMOUT,OPRINT,PRINT
11     COMMON,CONSTAN/PI,RADdeg,DEGRAD,RADsec,BOLTZ,SIN2I,ATCON
12     COMMON/LGIC/NIGHT,DAY,JLOCK,RESET,END
13     LOGICAL NIGHT, DAY, JLOCK, RESET, END, PRINT, SET
14     COMMON AL/ALGO/ZBC1,ZTOP,TURB,ZTRANS
15     DATA SET/,FALSE,/
16
17     IF(.NOT. SET) GO TO 1
18     SET=.FALSE.
19     DELT=1S
20     RESET=.FALSE.
21     TSEC=TIME*DELT
22     THRS=TSEC/3600.
23     IF(TINOUT.EQ.0.) GO TO 8
24     IF(THRS.LT. TINOUT) GOTO 10
25     IF(*4.5.EQ. TINOUT) GOTO 8
26     TSEC=1INOUT*3600
27     THREE=1MINT
28     DELT=1SEC-TIME
29     RESET=.TRUE.
30     PRINT=.TRUE.
31     TIMOUT=TINOUT+OPRINT
32     IF(.NOT. NIGHT) JLOCK=.TRUE.
33     GO TO 11
34
35     PRINT=.FALSE.
36     10 OLDMOD=AMOD(TIME,8.64E+04)
37
38     CHINU=ACOS(COSD(COS(TSEC*RADSEC)+SIND))
39
40     C DETERMINE WHETHER DAY, NIGHT, OR TWILIGHT.
41
42     C
43     IF(NIGHT .AND. CHINU .GT. SUNSET(K)) GO TO 15
44     IF(NIGHT .AND. CHINU .LT. SUNSET(K)) GO TO 5
45     IF(DAY .AND. CHINU .LT. SUNSET(11)) GO TO 13
46     IF(DAY .AND. CHINU .GT. SUNSET(11)) GO TO 9
47     IF(CHINU .GT. SUNSET(17)) NIGHT=.TRUE.
48     IF(CHINU .LT. SUNSET(17)) NIGHT=.FALSE.
49     IF(CHINU .EQ. SUNSET(17)) NIGHT=.TRUE.
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C FIND ALTITUDE OF DAY-NIGHT TRANSITION.
C
60      IF(DAY .OR. NIGHT) GO TO 26
      DO 6 J=1,K
      IF(CHINU .LE. SUNSET(J)) GO TO 7
      CONTINUE
65      7      ZTRANS=7/11
                  GO TO 26
C
C   RESET THE TIME STEP SO THAT THE NEXT TIME INCREMENT WILL BEGIN
C   EXACTLY AT SUNRISE AT THE UPPER BOUNDARY.
C
70      5      PELT=TRIGE-OLDMOD
                  TSEC=TIME+DELT
                  CHINU=SUNSET(K)
                  SET, TIME,
                  NIGHT=.FALSE.
                  GO TO 13
C
C   RESET THE TIME STEP SO THAT THE NEXT TIME INCREMENT WILL BEGIN
C   EXACTLY AT SUNSET AT THE LOWER BOUNDARY.
C
75      9      PELT=SET-OLDMOD
                  TSEC=TIME+DELT
                  CHINU=SUNSET(1)
                  SET=.TRUE.,
                  DAY=.FALSE.,
                  ZTRANS=50,
                  GO TO 26
                  13      15      ZTRANS=400,
                  26      CHI=CHINU
                  TIME=TSEC
                  RETURN
                  30      END
C
C   GIONEUT
1226    GIONEUT
1227    GIONEUT
1228    GIONEUT
1229    GIONEUT
1230    GIONEUT
1231    GIONEUT
1232    GIONEUT
1233    GIONEUT
1234    GIONEUT
1235    GIONEUT
1236    GIONEUT
1237    GIONEUT
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1248    GIONEUT
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1250    GIONEUT
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1252    GIONEUT
1253    GIONEUT
1254    GIONEUT
1255    GIONEUT
1256    GIONEUT
1257    GIONEUT
1258    GIONEUT
1259    GIONEUT

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SUBROUTINE CHEMION

C THIS SUBROUTINE COMPUTES THE SPECIES CONCENTRATIONS TO BE DETERMINED
C FROM CHEMISTRY ONLY. THE LARGEST NEGATIVE SPECIES IS COMPUTED
C FROM CHARGE BALANCE.

5 C

6 C DIMENSION H(1)

7 C

8 C COMMON NERS2,301,LC3N(1424,26),L(692,26),NZERO(692),UZERO(692)

9 C COMMON FORM(692),EWV(692),E(692),F(692),Z(692),Z(692),
2 C DECY(692),TEMP(692),STEMD(692),DETAZ(692),SUNSET(692)

10 C

11 C COMMON GRAV,SGND,TIME,X0,X1,RADIUS,LSPECI,KSPECI,NRFAC,K,
12 C X2,FWT,WT,IPAR,EXCNS,SCY1,SCY2,AREA2

13 C

14 C COMMON/LOGIC/NIGHT,DAY,JLOOK,RESET,END

15 C

16 C COMMON/RTIMES/DAYS,DELT,TMAX,TSET,TRISE

17 C

18 C COMMON/RAD,ECON,AM(88),BK(88),CK(215),CMB(6,215)

19 C

20 C COMMON/ATCONS/ATWT(26),MASS(26),SYMBOL(56),POLAR(26)

21 C

22 C LEVEL 3, OLDON,OLDON,OLDDU,RATES
REAL N NZERO
LOGICAL END

23 C EQUIVALENCE (U(1,21),H)

24 C COMPUTE OLD PROFILE.

25 C

26 C CALL MOYLEV(DION(1,27),NZERO,K)

27 DO 5 J=1,K
CALL CHEM(27,J)
N(I,27)=NZERO(J)+DELT*FORM(J)/(1.0+DELT*REMV(J))

28 5 CONTINUE

29 C

30 C COMPUTE ELECTRON, NEGATIVE ION AND CLUSTER ION PROFILES.

31 C

32 DO 45 J=1,K2
DO 10 I=1,3
CALL MOYLEV(OLDN(J,I+27),NZERO(I),1)

33 10 CONTINUE

34 DO 11 I=1,26
CALL MOYLEV(CION(J,I),NZERO(I+3),1)

35 11 CONTINUE

36 C COMPUTE ION, CLUSTER ION, ETC.

37 DO 45 I=1,24
DO 12 J=1,3
CALL MOYLEV(CION(J,I),NZERO(I+3),1)

38 12 CONTINUE

39 C COMPUTE TOTAL ION CONCENTRATION

40 C

41 C

42 C

43 C

44 C

45 C

46 C

47 C

48 C

49 C

50 C

51 C

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56 C

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      1001      E- 4      N-0.1881  *OC1
      1002      W-0.1778  *OC2
      1003      Z-0.2768  *OC3
      1004      GJD=47.8746
      1005      C     ALL NEGATIVE SPECIES.

      65          BIG=N(J,30)
      66          BIG=1
      67          DO 20 I=1,8
      68          IF(OCNC(J,I).LT.-BIG) GO TO 20
      69          BIG=CONC(J,I);
      70          JBIG=I+1
      71          CONTINUE
      72          C     SUM ALL POSITIVE SPECIES.
      73          C     SUMP=0.
      74          DO 30 I=1,6
      75          SUMP=SUMP+N(I,J,1)
      76          CONTINUE
      77          DO 31 I=23,29
      78          SUMP=SUMP+N(J,I)
      79          CONTINUE
      80          DO 35 I=9,24
      81          SUMP=SUMP+CCNC(J,I)
      82          CONTINUE
      83          DO 36 I=30,39
      84          SUMP=SUMP+CONC(J,I)
      85          CONTINUE
      86          C     SUM ALL NEGATIVE SPECIES EXCEPT LARGEST.
      87          C     SUMM=0.
      88          DO 40 I=1,9
      89          IF(I.EQ.JBIG) GO TO 40
      90          IF(I.EQ.1) SUMM=SUMM+N(J,30)
      91          IF(I.NE.1) SUMM=SUMM+CONC(J,I-1)
      92          CONTINUE
      93          DEN=SUMP-SUMM
      94          IF(DEN.GT.0.) GO TO 79
      95          WRITE(6,75) Z(J),L,JBIG,SUMP,SUMM,DELT
      96          WRITE(6,76) N(J,30),(CONC(J,I),I=1,8)
      97          FORMAT(1P9E13.5)
      98          FORMAT(1* ERROR IN CHEMION. ALI=*,FB.2,14,13,1P3E12.5)
      99          END* .TRUE.
      100         RETURN
      101        IF(JBIG.EQ.-1) N(J,20)=DENS
      102        IF(JBIG.NE.-1) CONC(J,JBIG)=DENS
      103        DO 65 I=1,29
      104        IF(I.EQ.3) DFNS=N(J,I+27)
      105        IF(I.EQ.3) DFNS=CONC(J,I-3)
      106        IF(DENS.LT.0.) I=J GO TO 65
      107        IF(H(I).EQ.0.) GO TO 65
      108        IF((ABS(1./H(I))-DENS)/H(I)).GT.0.20) GO TO 60
      109        GO TO 45
      110        CONTINUE
      111        N=0
      112        CONTINUE
      113        WRITE(6,*) ' CALCULATIONS NOT CONVERGING IN CHEMION. '
      114        WRITE(6,*) ' TRYING ANOTHER CONVERGENCE TEST'
      115        GO TO 10
      116        END

```

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115      81   FORMAT(• FAILURE AT ALTITUDE•,OPF8.2,• FOR SPECIES•,A10,1P2E13.5) GIONEUT
          END•.TRUE.
          RETURN
        45   CONTINUE
C
120      C COMPUTE THE MAJOR POSITIVE IONS AND ELECTRONS ABOVE 120 KM.
C
        KB=K2+1
        CALL MOVELEV(OLDN(1,28),E,K)
        CALL MOVELEV(OLDN(1,29),F,K)
        DO 105 J=KB,K
        DO 100 L=1,500
        DO 110 I=1,2
        H(I)=N(J,I+27)
        CALL CHEM(I+27,J)
        IF(I .EQ. 1) DENS=E(J)
        IF(I .EQ. 2) DENS=F(J)
        H(I+3)=DENS+DELT*FORM(J)/(1.0+DELT*REMV(J))
        110  CONTINUE
        DO 91 I=1,2
        N(J,I+27)=H(I+3)
        91   CONTINUE
        DO 95 I=1,2
        IF(H(I) .EQ. 0.) GO TO 95
        IF(ABS(1.0-(N(J,I+27)/H(I))) .GT. 0.1) GO TO 100
        95   CONTINUE
        DO 55 J=KB,K
        SUMP=0.
        DO 50 I=1,6
        SUMP=SUMP+N(J,I)
        50   CONTINUE
        DO 51 I=28,29
        SUMP=SUMP+N(J,I)
        51   CONTINUE
        N(J,30)=SUMP
        55   CONTINUE
        RETURN
        END
135
140
145
150
155

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      1415
      GIONEUT 1415
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      GIONEUT 1469
      GIONEUT 1470
      GIONEUT 1471

10    G  COMMON FORM(692),REW(692),E(692),Z(692),
      G  E(692),Z(692),DELTA(692),DELTA(692),SUNSE1(692)
      G  S  COMMON GRAV,LSD,SIND,TIME,CAT,RADIUS,ISPECI,KSPECI,NREAC,KTOP,
      G  F2,FMIN,FLFS,CX,NOON,SCXI,NREAC2
      G  COMMON RATECON,AK(88),BK(88),CK(88),DK(215),CYMB(6,215)
      G  COMMON,FSCN(692,30),LOCCON(424,26),FLDU(692,26),RATES(88,692)
      G  C  COMMON,FCN,NIGHT,DAY,FLDOOR,RESET,END
      G  C  C 1000,1000,LOCCON,O-DU,RATES
      G  C  LOGICAL,NIGHT
      H  ALT M N
      G  DO 500 1=1,1,ISPECI
      G  IF(I .EQ. 1) GO TO 200
      G  Y(I)=H(ALT,I)
      G  GO TO 500
      G  200  IF((ALT .GT. K2) GO TO 475
      G  T1=FCNC(-ALT,1-30)
      G  GO TO 500
      G  475  Y(1)=C,
      G  CONTINUE
      G  Y(57)=0,
      G  DO 600 1=1,1,ISPECI
      G  Y(57)+Y(57)+Y(1)
      G  600  CONTINUE
      G  DO 205 1=1,NREAC
      G  K(1)=DK(1)
      G  205  CONTINUE
      G  IF(NOT. NIGHT) GO TO 210
      G  K(96)=K(59)*K(104)=K(105)*K(110)*C
      G  K(113)=K(114)*K(118)*K(120)*K(123)*C
      G  CALL MOVEL(V(RATEFS(1,1)AT),K(128),B,B
      G  IF(TALI LE. K2) GC TO 215
      G  K(31)=K(32)*K(35)*C
      G  K(16)=K(16)*K(152)*K(165)*K(166)*C
      G  110  0(1,2,3,4,5,6,7,B,S,10,11,12,13,14,15,16,17,18,19,20,21,22,23
      G  ,4,25,26,27,28,29,30,31,32,33,34,35,36,37,38,39,40,41,42,43,44,45,
      G  ,45,46,47,48,49,50,51,52,53,54,55,56),KIND
      G  1247,1248,1249
      G  1249,1250,1251,1252,1253,1254,1255,1256,1257,1258,1259,1260,1261,1262,1263,1264,1265,1266,1267,1268,1269,1270,1271

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RD-A157 122

A COMPUTER CODE FOR A ONE-DIMENSIONAL DYNAMIC MODEL OF
THE MESOSPHERE AND... (U) AIR FORCE GEOPHYSICS LAB
HANSOM AFB MA T J KENESHEA ET AL. 07 MAR 84

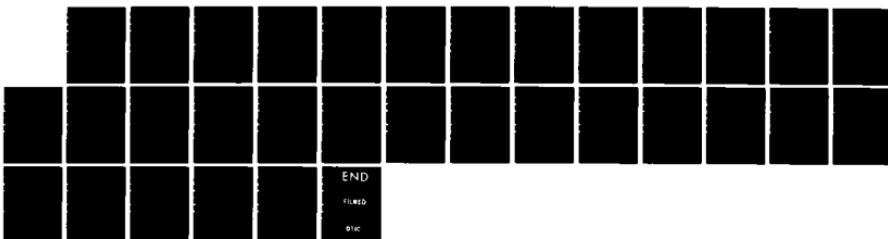
3/3

UNCLASSIFIED

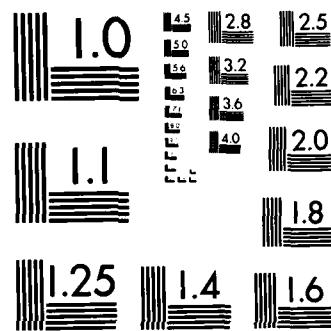
AFGL-TR-84-0183

F/G 4/1

NL



END
FILED
SIC



MICROCOPY RESOLUTION TEST CHART
STANDARDS-1963-A

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RMOV=(Y(8)*(+K(45)+K(46))+Y(23)*(+K(47)+K(48))+K(49)*Y(18)) GIONEUT 1472
FARM=(+K(191)*Y(21)) GIONEUT 1473
GO TO 100 GIONEUT 1474
3 CONTINUE GIONEUT 1475
RMOV=(+K(40)*Y(11)+K(127)*Y(25)+K(153)*Y(30)+K(168)*Y(23)+K(169)*Y GIONEUT 1476
1(B) FARM=(Y(7)*(+K(41)*Y(1)+K(185)+K(211))+Y(8)*(+K(45)*Y(2)+K(209))) GIONEUT 1477
FARM=(Y(8)*(+K(43)+K(44)+K(46)*Y(2)+K(126)*Y(29)+K(167)*Y(29)+K(209))) GIONEUT 1478
GO TO 100 GIONEUT 1479
4 CONTINUE GIONEUT 1480
RMOV=(+K(33)*Y(18)+K(34)*Y(16)+K(35)*Y(19)+K(152)*Y(30)+K(162)*Y(8) GIONEUT 1481
1)*Y(57)) FARM=(Y(8)*(+K(43)+K(44)+K(46)*Y(2)+K(126)*Y(29)+K(167)*Y(29)+K(209))) GIONEUT 1482
FARM=(Y(8)*(+K(169)*Y(3)+K(186)+K(214)+Y(39)*(+K(36)+Y(7)+K(163)*Y(57)+K(16 GIONEUT 1483
18)+K(169)*Y(3)+K(186)+K(214)+Y(39)*(+K(36)+Y(7)+K(163)*Y(57)+K(16 GIONEUT 1484
14)*Y(20))+K(127)*Y(3)*Y(25)+K(207)*Y(20))) GIONEUT 1485
GO TO 100 GIONEUT 1486
5 CONTINUE GIONEUT 1487
RMOV=(Y(30)*(+K(150)+K(151))+K(165)*Y(25)*Y(23)+K(166)*Y(13)*Y(57)) GIONEUT 1488
1) FARM=(Y(18)*(+K(33)*Y(4)+K(50)*Y(40)+K(188)*Y(7)*(+K(38)*Y(28)+K( GIONEUT 1489
139)*Y(28))+K(34)*Y(4)*Y(16)+K(42)*Y(6)*Y(8)+K(168)*Y(3)*Y(23))) GIONEUT 1490
GO TO 100 GIONEUT 1491
6 CONTINUE GIONEUT 1492
RMOV=(Y(8)*{+K(42)+K(43)+K(44)+K(154)*Y(30)}) GIONEUT 1493
FARM=(Y(2)*(+K(47)*Y(23)+K(49)*Y(18))+K(189)*Y(16)+K(208)*Y(23)) GIONEUT 1494
GO TO 100 GIONEUT 1495
7 CONTINUE GIONEUT 1496
RMOV=(Y(57)*(Y(7)*(+K(128)+K(129)*Y(8)+K(139)*Y(16)+K(143) GIONEUT 1497
1*Y(18)+K(148)*Y(26))+Y(14)*(+K(132)+K(133))+Y(28)*{+K(38)+K(39)}*Y GIONEUT 1498
1(32)*(+K(100)+K(101))+Y(34)*(+K(107)+K(108))+K(2)*Y(10)+K(3)*Y(12) GIONEUT 1499
1+K(4)*Y(19)+K(20)*Y(55)+K(31)*Y(30)+K(36)*Y(39)+K(41)*Y(1)+K(97)*Y GIONEUT 1500
1(31)+K(106)*Y(33)+K(111)*Y(37)+K(115)*Y(38)+K(121)*Y(36)+K(130)*Y( GIONEUT 1501
19)+K(131)*Y(15)+K(185)+K(206)+K(211)+K(212)) GIONEUT 1502
FARM=Y(8)*(+K(14)+Y(17)+K(42)*Y(6)+K(45)*Y(2)+K(126)*Y(29)+K(137) GIONEUT 1503
1*Y(16)+K(169)*Y(3)+K(192)+K(193)+K(209)+K(210))+Y(30)*(Y(4) GIONEUT 1504
1*(+K(152)+K(152))+Y(5)*(+K(150)+K(151))+K(153)*Y(3)+K(156)*Y(40)+ GIONEUT 1505
1Y(16)*+K(9)*Y(19)+K(34)*Y(4)+K(138)*Y(18))+Y(9)*(+K(98)*Y(31)+K(1 GIONEUT 1506
194))+Y(18)*{+K(49)*Y(2)+K(199)+K(122)*Y(10)+Y(10)+K(16)*Y(27)*Y(57) GIONEUT 1507
1)+K(40)*Y(3)*Y(11)+K(96)*Y(31)+K(120)*Y(36)+K(125)*Y(29)+K(12 GIONEUT 1508
104)*Y(25)}) GIONEUT 1509
GO TO 100 GIONEUT 1510
8 CONTINUE GIONEUT 1511
RMOV=(Y(8)*(Y(30)*(+K(160)+K(160)+Y(31)*(+K(179)+K(179)*Y(32)*{+ K(180)+K(180)+Y(63)*Y(33)+Y(39)+K(114)*Y(38)-K(205))) GIONEUT 1512
12)*Y(10)+K(14)*Y(6)*{+K(42)+K(43)+K(44)+Y(57)*(+K(129)*Y(7)+K(13 GIONEUT 1513
16)*Y(11)+K(162)*Y(4)+Y(2)*(+K(45)+K(46))+K(14)*Y(17)+K(32)*Y(23)* GIONEUT 1514
1Y(30)+K(126)*Y(29)+K(137)*Y(16)+K(167)*Y(28)+K(169)*Y(3)+K(186)+K( GIONEUT 1515
1Y(38)+K(193)+K(209)+K(210)+K(214)) GIONEUT 1516
FARM=(Y(7)*(Y(34)*(+K(107)+K(108)+Y(9)*(+K(130)+K(130))+K(1517) GIONEUT 1517
12)*Y(10)+K(3)*Y(12)+K(4)*Y(19)+K(97)*Y(31)+K(101)*Y(32)+K(106)*Y(3 GIONEUT 1518
13)*K(115)*Y(38)+K(121)*Y(36)+K(128)*Y(7)*Y(57)*(+K(129)*Y(14)+Y(9)* GIONEUT 1519
1Y(12)*(+K(145))*K(7)*Y(11)+K(8)*Y(16)+K(16)+K(102)*Y(32)+K(116)* GIONEUT 1520
1Y(38)+K(119)*Y(.35)+K(124)*Y(31)+K(140)*Y(10)+K(142)*Y(18)+K(161)*Y GIONEUT 1521
1*(30)+K(194)*Y(20)*{+K(103)+K(103)+Y(39)+(Y(39)*(+K(164)+K(164))+Y(9) GIONEUT 1522
1K(1B)*Y(57)+K(19)+Y(12)*{+K(5)*Y(11)+K(13)*Y(10)+K(144)*Y(12)+K(1 GIONEUT 1523
198))+Y(8)*Y(8)*{+K(160)*Y(30)+K(179)*Y(31)+K(180)*Y(32)})+Y(13)*Y(32) GIONEUT 1524
1Y(50)*{+K(53)+K(54)+K(37)*Y(39)+Y(4)*(+K(33)*Y(18)+K(35)*Y(19)+Y(14) GIONEUT 1525
1Y(25)*(-K(1C9)*Y(34)-K(182)*Y(33)+Y(57)*(+K(163)*Y(39)+K(63)*Y(3) GIONEUT 1526
14))+K(111)*Y(16)-Y(19)+K(99)*Y(32)+K(105)*Y(33)+K(114)*Y(38)-K(205)) GIONEUT 1527
1528

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115      1*Y(56)
      GO TO 100
9      CONTINUE
      RMDV=(Y(31)*(+K(98)+K(124))+K(7)*Y(11)+K(B)*Y(16)+K(102)*Y(32)+K(1
      116)*Y(38)+K(119)*Y(35)+K(130)*Y(7)+K(140)*Y(10)+K(142)*Y(18)+K(145)
      1)*Y(12)+K(161)*Y(30)+K(194+K(195))
      FARM=(Y(7)*(+K(36)*Y(39)+K(100)*Y(32)+K(129)*Y(8)*Y(57))+K(104)*Y(
      133))
      GO TO 100
10     CONTINUE
      RMDV=(Y(10)*(+K(12)+K(2)*Y(7)+K(13)*Y(12)+K(21)*Y(55)+K(140
      1)*Y(9)+K(141)*Y(14)+K(147)*Y(26))
      FARM=(Y(37)*(+K(64)*Y(41)+K(68)*Y(42)+K(72)*Y(43)+K(76)*Y(44)+K(80
      1)*Y(45)+K(92)*Y(48)+K(94)*Y(49)+Y(30)*(+K(24)*Y(41)+K(25)*Y(42)+K
      1)(26)*Y(43)+K(44)+K(28)*Y(45)+K(29)*Y(47)+Y(11)*Y(12)+(+K(6
      1)+K(6)*K(7)*Y(9)+K(135)*Y(14)+K(146)*Y(19)+Y(13)*(Y(27)*(+K(15)+
      1K(15))+K(54)*Y(50)+K(61)*Y(47)+K(196)+Y(36)*(+K(63)*Y(41)+K(67)*Y
      1(42)+K(71)*Y(43)+K(75)*Y(44)+K(79)*Y(45)+Y(7)*(+K(3)*Y(12)+K(20)*
      1Y(55)+K(131)*Y(15)+K(132)*Y(14)+Y(14)*(+K(197)+K(197))+K(17)*Y(27
      1)*Y(15)+K(145)*Y(12)+Y(9)+K(202)*Y(55))
      GO TO 100
11     CONTINUE
      RMQV=(Y(12)*(+K(5)+K(6))+Y(14)*(+K(134)+K(135))+K(7)*Y(9)+K(40)*Y(
      13)+K(136)*Y(8)*Y(57)+K(146)*Y(19)+K(190))
      FARM=(Y(35)*(+K(62)*Y(41)+K(66)*Y(42)+K(70)*Y(43)+K(74)*Y(44)+K(78
      1)*Y(45))+Y(7)*(+K(2)*Y(10)+K(41)*Y(11)+K(131)*Y(15))+Y(30)*(+K(29)*
      1Y(47)+K(30)*Y(49)+K(155)*Y(11)+K(17)*Y(27)*Y(15)+K(147)*Y(17)
      1)+K(196)*Y(13)+K(198)*Y(12))
      GO TO 100
12     CONTINUE
      RMDV=(Y(11)*(+K(5)+K(6))+Y(12)*(+K(144)+K(144))+K(3)*Y(7)+K(13)*Y(
      110)+K(145)*Y(9)+K(198))
      FARM=(Y(38)*(+K(69)*Y(42)+K(42)*Y(73)*Y(43)+K(77)*Y(44)+K(B1)*Y(45)+K(93
      1)*Y(48)+K(95)*Y(49)+Y(14)*Y(14)+K(132)*Y(7)+K(134)*Y(11)+K(141)*Y(10)
      1)+K(65)*Y(42)*Y(32)+K(136)*Y(11)*Y(8)*Y(57)+K(140)*Y(10)*Y(9))
      GO TO 100
13     CONTINUE
      RMDV=(Y(57)*(+K(52)*Y(44)+K(166)*Y(5)+K(170)*Y(41)+K(172)*Y(42)+K(
      1173)*Y(43)+K(177)*Y(51)+K(178)*Y(52)+Y(48)*(+K(58)*Y(23)+K(60))+Y
      1(50)*Y(53)+K(54)+K(15)*Y(27)+K(37)*Y(39)+K(55)*Y(53)+K(56)*Y(54
      1)+K(59)*Y(49)+K(61)*Y(47)+K(196))
      FARM=(Y(30)*(Y(52)*(+K(158)+K(158))+K(25)*Y(42)+K(29)+K(157)
      1*Y(51)+Y(10)*(+K(12)*Y(10)+K(13)*Y(12)+K(21)*Y(55)+K(141)*Y(14))+
      1Y(37)*(+K(64)*Y(41)+K(84)*Y(51)+K(92)*Y(48)+Y(57)*(+K(174)*Y(43)+
      1K(175)*Y(44)+K(176)*Y(45)+Y(14)*Y(14)*(+K(133)*Y(7)+K(135)*Y(11)+Y(32)
      1*(+K(65)*Y(42)+K(65)*Y(42)+Y(35)*(+K(62)*Y(41)+K(82)*Y(51))+Y(36)
      1*(+K(63)*Y(41)+K(83)*Y(51)+K(93)*Y(48)*Y(38)))
      FARM=FARM+2.*((Y(42)*(K(66)*Y(35)+K(67)*Y(36)+K(68)*Y(37)
      1+K(69)*Y(38))+Y(49)*(K(30)*Y(30)+K(94)*Y(37)+K(95)*Y
      2*(38)+Y(52)*(K(85)*Y(35)+K(86)*Y(36)+K(87)*Y(37))
      3+(K(26)*Y(43)*Y(30))-3.0*(Y(43)*K(70)*Y(35)+K(71)*
      4Y(36)+K(72)*Y(37)+K(73)*Y(38))+Y(53)*(K(88)*Y(35)
      5+K(89)*Y(36)+K(90)*Y(37)+K(159)*Y(30)+K(27)*Y(44)
      6*Y(30)+4.0*(Y(44)*(K(74)*Y(35)+K(75)*Y(36)+K(76)*Y(
      737)+K(77)*Y(39)+K(28)*Y(45)*Y(30)+5.0*(Y(45)*
      6K(78)*Y(35)+K(79)*Y(36)+K(60)*Y(37)+K(81)*Y(38)))
      GO TO 100

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14 CONTINUE
    RMOV=(Y(7)*(+K(132)+K(133))+Y(11)*(+K(134)+K(135)))*K(141)*Y(10)+K(
1197)
175
    FARM=(+K(91)*Y(48)*Y(36)+K(144)*Y(12))*Y(12))
    GO TO 100
15 CONTINUE
    RMOV=(+K(17)*Y(27)+K(131)*Y(7))
    FARM=(Y(30)*(+K(24)*Y(41)+K(25)*Y(42)+K(26)*Y(43)+K(27)*Y(44)+K(28)
1)*Y(45))+Y(11)*(+K(5)*Y(12)+K(134)-Y(14)))
    GO TO 100
16 CONTINUE
    RMOV=(Y(19)*(+K(10)+K(11))+K(8)*Y(9)+K(34)*Y(4)+K(137)*Y(8))-K
1*(138)*Y(18)+K(139)*Y(7)*Y(57)+K(189)
    FARM=(Y(30)*(Y(28)*(+K(22)+K(22)+K(149)*Y(17)+K(150)*Y(5)+K(154)*
1*Y(6))+Y(23)*(+K(47)*Y(2)+K(168)*Y(3)+K(208)+K(215))+Y(7)*(+K(1)*Y(
117)+K(38)*Y(28)+K(43)*Y(6)*Y(8)+K(199)*Y(18)))
    GO TO 100
17 CONTINUE
    RMOV=(+K(1)*Y(7)+K(14)*Y(8)+K(149)*Y(30))
    FARM=(Y(28)*(Y(30)*(+K(23)+K(23))+K(39)*Y(7))+K(44)*Y(6)+K(B)+K(15
11)*Y(5)*Y(30)+K(215)*Y(23))
    GO TO 100
18 CONTINUE
    RMOV=(+K(33)*Y(4)+K(49)*Y(2)+K(50)*Y(40)+K(117)*Y(38)+K(122)*Y(56)
1+K(138)*Y(16)+K(142)*Y(9)+K(143)-Y(7)*Y(57)+K(184)*Y(37)+K(188)+K(
1199))
    FARM=(Y(16)*(Y(19)*(+K(10)+K(10))+K(8)*Y(9)+K(137)*Y(8)+K(139)*Y(7
1)*Y(57))+Y(30)*(+K(156)*Y(40)+K(157)*Y(51)+K(158)*Y(52)+K(159)*Y(5
13)+Y(19)*(+K(4)*Y(7)+K(146)*Y(11)+K(200))+Y(35)*(+K(82)*Y(51)+K(8
15)*Y(52)+K(88)*Y(53))+K(14)*Y(17)-Y(8)+K(202)*Y(55)+K(205)*Y(56)))
    GO TO 100
19 CONTINUE
    RMOV=(Y(16)*(+K(9)+K(10)+K(11))+K(4)*Y(7)+K(35)*Y(4)+K(112)*Y(37)+
1K(146)*Y(11)+K(200))
    FARM=(Y(36)*(Y(51)*(+K(83)+K(83))+Y(52)*(+K(86)+K(86))+Y(53)*(+K(8
19)+K(89))+K(63)*Y(41)+K(67)*Y(42)+K(71)*Y(43)+K(75)*Y(44)+K(79)*Y(
145)+Y(35)*(+K(62)*Y(41)+K(66)*Y(42)+K(70)*Y(43)+K(74)*Y(44)+K(78))
1*Y(45)+K(82)*Y(51)+K(85)*Y(52)+K(88)*Y(53)+K(118))+Y(18)*(Y(56)+Y(57)
1K(122)+K(122))+K(50)*Y(40)+K(142)*Y(9)+K(143)*Y(7)+Y(37)*(+K(84)*Y(51)+K(87)*Y(52)+K(90)*Y(53))+Y(55)*(+K(20)*Y(7)+K(21))-Y(10
1)))
    GO TO 100
20 CONTINUE
    RMOV=(+K(18)*Y(57)+K(19)+K(103)+Y(32)+K(164)*Y(39)+K(207))
    FARM=(+K(195)*Y(9))
    GO TO 100
21 CONTINUE
    RMOV=(+K(191))
    FARM=(Y(2)*(Y(8)*(+K(45)+K(46))+Y(23)*(+K(47)+K(48))+K(49)*Y(18)))
    GO TO 100
22 CONTINUE
    RMOV=0.
    FARM=0.
    GO TO 100
23 CONTINUE
    RMOV=(Y(2)*(+K(47)+K(48))+Y(23))-(-K(171)-Y(41)+K(171)*Y(215))
1*Y(29)+K(168)*Y(3)+K(187)+K(208)+K(213)+K(215))

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230      FARM=(Y(16)*(+K(11)*Y(19)+K(138)*Y(18))+K(57)*Y(46)*Y(25)+K(167)*Y
1*(28)*Y(8)+K(171)*Y(41)*Y(23)*Y(201)+K(24)) GIONEUT 1643
      GO TO 100 GIONEUT 1644
24      CONTINUE GIONEUT 1645
      RMDV=(+K(201))
      FARM=(+K(9)*Y(16)*Y(19))
      GO TO 100 GIONEUT 1646
25      CONTINUE GIONEUT 1647
      RMDV=(+K(51)*Y(42)*Y(57)+K(57)*Y(46)+K(109)*Y(34)+K(127)*Y(3)+K(16
15)*Y(5)*Y(23)+K(181)*Y(32)*Y(8)+K(182)*Y(33)+K(203)+K(204)) GIONEUT 1650
      FARM=(Y(37)*(+K(64)*Y(41)+K(68)*Y(42)+K(72)*Y(43)+K(76)*Y(44)+K(80
1)*Y(45)+K(84)*Y(51)+K(87)*Y(52)+K(90)*Y(53)+K(110)+K(111)*Y(17)+K(1
112)*Y(19)+K(184)*Y(18)+Y(38)*(+K(69)*Y(42)+K(73)*Y(43)+K(77)*Y(44
1)+K(81)*Y(45)+K(113)+K(114)+K(116)*Y(9)+K(117)*Y(18))+Y(13)*(+K(56
1)*Y(54)+K(59)*Y(49)+K(60)*Y(48))+Y(26)*(+K(147)*Y(10)+K(148)*Y(7)*
1Y(57)+K(30)*Y(45)*Y(30)+K(91)*Y(48)*Y(36)) GIONEUT 1655
      FARM=FARM+2.0*(Y(37)*(K(92)*Y(48)+K(54)*Y(49))+Y(38)*(1K(93)*Y(48)+K(95)*Y(49)))
      GO TO 100 GIONEUT 1656
26      CONTINUE GIONEUT 1657
      RMDV=(+K(147)*Y(10)+K(148)*Y(7)*Y(57))
      FARM=(Y(25)*(+K(127)*Y(3)+K(203)+K(204)))
      GO TO 100 GIONEUT 1658
27      CONTINUE GIONEUT 1659
      RMDV=(+K(15)*Y(13)+K(16)*Y(57)+K(17)*Y(15))
      FARM=(+K(193)*Y(8)+K(195)*Y(9)+K(200)*Y(19)+K(201)*Y(24)+K(203)*Y(
125))
      GO TO 100 GIONEUT 1660
28      CONTINUE GIONEUT 1661
      RMDV=(Y(7)*(+K(38)+K(39))+Y(30)*(+K(22)+K(23))+K(167)*Y(8))
      FARM=(Y(23)*(+K(48)*Y(2)+K(125)*Y(29)+K(187)+K(213)))
      GO TO 100 GIONEUT 1662
29      CONTINUE GIONEUT 1663
      RMDV=(+K(125)*Y(23)+K(126)*Y(8))
      FARM=(Y(7)*(+K(206)+K(212))+K(210)*Y(8))
      GO TO 100 GIONEUT 1664
30      CONTINUE GIONEUT 1665
      RMDV=(Y(5)*(+K(150)+K(151))+Y(8)*(+K(32)+K(160)*Y(8))+Y(28)*
1(+K(22)+K(23))+K(24)*Y(41)+K(25)*Y(42)+K(26)*Y(43)+K(27)*Y(44)+K(2
18)*Y(45)+K(29)*Y(47)+K(30)*Y(49)+K(31)*Y(7)+K(152)*Y(4)+K(153)*Y(3
1)+K(154)*Y(6)+K(155)*Y(1)+K(156)*Y(40)+K(157)*Y(51)+K(158)*Y(52)+K
1(159)*Y(53)+K(161)*Y(9)) GIONEUT 1666
      FARM=(Y(7)*(+K(97)+K(100)+Y(31)+K(185)+K(206))+Y(20)*(+K(103
1)*Y(32)+K(207))+Y(23)*(K(187)+K(96)*Y(31)+K(99)*Y(32)+K(1
104)*Y(33)+K(114)*Y(38)+K(118)*Y(35)+K(123)*Y(36)+K(186)*Y(8)+K(188
1)*Y(18)+K(189)*Y(16)+K(190)*Y(11)+K(191)*Y(21))
      FARM=FARM+Y(23)*(K(208)+K(213))+Y(8)*(K(209)+K(210)+K(214))
      GO TO 100 GIONEUT 1667
31      CONTINUE GIONEUT 1668
      RMDV=(Y(9)*(+K(98)+K(124))+K(96)+K(97)*Y(7)+K(179)*Y(8))
      FARM=(Y(7)*(+K(31)*Y(30)+K(101)*Y(32)+K(108)*Y(34)+K(105)*Y(33)+K
1(110)*Y(37)+K(161)*Y(39)+Y(30))
      GO TO 100 GIONEUT 1669
32      CONTINUE GIONEUT 1670
      RMDV=(Y(7)*(+K(100)+K(101))+Y(8)*(+K(180)*Y(8)+K(181)*Y(25))+K(65)
1*Y(42)+K(99)+K(102)*Y(9)+K(103)*Y(20)) GIONEUT 1671
      FARM=(Y(7)*(+K(31)*Y(30)+K(101)*Y(32)+K(108)*Y(34)+K(105)*Y(33)+K
1(110)*Y(37)+K(161)*Y(39)+Y(30))
      GO TO 100 GIONEUT 1672
33      CONTINUE GIONEUT 1673
      RMDV=(Y(7)*(+K(100)+K(101))+Y(8)*(+K(180)*Y(8)+K(181)*Y(25))+K(65)
1*Y(42)+K(99)+K(102)*Y(9)+K(103)*Y(20)) GIONEUT 1674
      FARM=(Y(7)*(+K(31)*Y(30)+K(101)*Y(32)+K(108)*Y(34)+K(105)*Y(33)+K
1(110)*Y(37)+K(161)*Y(39)+Y(30))
      GO TO 100 GIONEUT 1675
34      CONTINUE GIONEUT 1676
      RMDV=(Y(7)*(+K(100)+K(101))+Y(8)*(+K(180)*Y(8)+K(181)*Y(25))+K(65)
1*Y(42)+K(99)+K(102)*Y(9)+K(103)*Y(20)) GIONEUT 1677
      FARM=(Y(7)*(+K(31)*Y(30)+K(101)*Y(32)+K(108)*Y(34)+K(105)*Y(33)+K
1(110)*Y(37)+K(161)*Y(39)+Y(30))
      GO TO 100 GIONEUT 1678
35      CONTINUE GIONEUT 1679
      RMDV=(Y(7)*(+K(100)+K(101))+Y(8)*(+K(180)*Y(8)+K(181)*Y(25))+K(65)
1*Y(42)+K(99)+K(102)*Y(9)+K(103)*Y(20)) GIONEUT 1680
      FARM=(Y(7)*(+K(31)*Y(30)+K(101)*Y(32)+K(108)*Y(34)+K(105)*Y(33)+K
1(110)*Y(37)+K(161)*Y(39)+Y(30))
      GO TO 100 GIONEUT 1681
36      CONTINUE GIONEUT 1682
      RMDV=(Y(7)*(+K(100)+K(101))+Y(8)*(+K(180)*Y(8)+K(181)*Y(25))+K(65)
1*Y(42)+K(99)+K(102)*Y(9)+K(103)*Y(20)) GIONEUT 1683
      FARM=(Y(7)*(+K(31)*Y(30)+K(101)*Y(32)+K(108)*Y(34)+K(105)*Y(33)+K
1(110)*Y(37)+K(161)*Y(39)+Y(30))
      GO TO 100 GIONEUT 1684
37      CONTINUE GIONEUT 1685
      RMDV=(Y(7)*(+K(100)+K(101))+Y(8)*(+K(180)*Y(8)+K(181)*Y(25))+K(65)
1*Y(42)+K(99)+K(102)*Y(9)+K(103)*Y(20)) GIONEUT 1686
      FARM=(Y(7)*(+K(31)*Y(30)+K(101)*Y(32)+K(108)*Y(34)+K(105)*Y(33)+K
1(110)*Y(37)+K(161)*Y(39)+Y(30))
      GO TO 100 GIONEUT 1687
38      CONTINUE GIONEUT 1688
      RMDV=(Y(7)*(+K(100)+K(101))+Y(8)*(+K(180)*Y(8)+K(181)*Y(25))+K(65)
1*Y(42)+K(99)+K(102)*Y(9)+K(103)*Y(20)) GIONEUT 1689
      FARM=(Y(7)*(+K(31)*Y(30)+K(101)*Y(32)+K(108)*Y(34)+K(105)*Y(33)+K
1(110)*Y(37)+K(161)*Y(39)+Y(30))
      GO TO 100 GIONEUT 1690
39      CONTINUE GIONEUT 1691
      RMDV=(Y(7)*(+K(100)+K(101))+Y(8)*(+K(180)*Y(8)+K(181)*Y(25))+K(65)
1*Y(42)+K(99)+K(102)*Y(9)+K(103)*Y(20)) GIONEUT 1692
      FARM=(Y(7)*(+K(31)*Y(30)+K(101)*Y(32)+K(108)*Y(34)+K(105)*Y(33)+K
1(110)*Y(37)+K(161)*Y(39)+Y(30))
      GO TO 100 GIONEUT 1693
40      CONTINUE GIONEUT 1694
      RMDV=(Y(7)*(+K(100)+K(101))+Y(8)*(+K(180)*Y(8)+K(181)*Y(25))+K(65)
1*Y(42)+K(99)+K(102)*Y(9)+K(103)*Y(20)) GIONEUT 1695
      FARM=(Y(7)*(+K(31)*Y(30)+K(101)*Y(32)+K(108)*Y(34)+K(105)*Y(33)+K
1(110)*Y(37)+K(161)*Y(39)+Y(30))
      GO TO 100 GIONEUT 1696
41      CONTINUE GIONEUT 1697
      RMDV=(Y(7)*(+K(100)+K(101))+Y(8)*(+K(180)*Y(8)+K(181)*Y(25))+K(65)
1*Y(42)+K(99)+K(102)*Y(9)+K(103)*Y(20)) GIONEUT 1698
      FARM=(Y(7)*(+K(31)*Y(30)+K(101)*Y(32)+K(108)*Y(34)+K(105)*Y(33)+K
1(110)*Y(37)+K(161)*Y(39)+Y(30))
      GO TO 100 GIONEUT 1699

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FARM=(Y(7)*(+K(106)*Y(33)+K(111)*Y(37))+Y(8)*(Y(30)*(+K(32)*Y(23))+  
1K(160)*Y(8)))+K(113)*Y(38)+K(124)*Y(31)+Y(9)+K(183)+Y(57)) * Y(57) +  
GIONEUT 1700  
GO TO 100  
33 CONTINUE  
RMOV=(+K(104)+K(05)+K(106)*Y(7)+K(182)*Y(25))  
FARM=(Y(9)*(+K(98)*Y(31)+K(102)*Y(32)+K(116)*Y(38))+K(107)*Y(34))*Y  
1(7)+K(179)*Y(31)*Y(8)*Y(8))  
GO TO 100  
34 CONTINUE  
RMOV=(Y(7)*(+K(107)+K(108))+K(109)*Y(25)+K(183)*Y(57))  
FARM=(+K(180)*Y(32)+Y(8)*Y(8))  
GO TO 100  
35 CONTINUE  
RMOV=(+K(62)*Y(41)+K(66)*Y(42)+K(70)*Y(43)+K(74)*Y(44)+K(78)*Y(45)  
1+K(B2)*Y(51)+K(85)*Y(52)+K(88)*Y(53)+K(118)*Y(59))  
FARM=(Y(36)*(+K(120)+K(121)*Y(7))+K(184)*Y(37)*Y(18))  
GO TO 100  
36 CONTINUE  
RMOV=(+K(63)*Y(41)+K(67)*Y(42)+K(71)*Y(43)+K(75)*Y(44)+K(79)*Y(45)  
1+K(83)*Y(51)+K(86)*Y(52)+K(89)*Y(53)+K(91)*Y(48)+K(120)+K(121)*Y(7)  
1)+K(123))  
FARM=(+K(112)*Y(37)+Y(19)+K(117)*Y(38)+Y(18)+K(119)*Y(35)+Y(9))  
GO TO 100  
37 CONTINUE  
RMOV=(+K(64)*Y(41)+K(68)*Y(42)+K(72)*Y(43)+K(76)*Y(44)+K(80)*Y(45)  
1+K(84)*Y(51)+K(87)*Y(52)+K(90)*Y(53)+K(92)*Y(48)+K(94)*Y(49)+K(110)  
1)+K(111)*Y(7)+K(112)*Y(19)+K(184)*Y(18))  
FARM=(+K(115)*Y(38)+Y(7)+K(182)+Y(33)+Y(25))  
GO TO 100  
38 CONTINUE  
RMOV=(+K(69)*Y(42)+K(73)*Y(43)+K(77)*Y(44)+K(81)*Y(45)+K(93)*Y(48)  
1+K(95)*Y(49)+K(113)+K(114)+K(115)*Y(7)+K(116)*Y(9)+K(117)*Y(18))  
FARM=(Y(25)*(+K(109)*Y(34)+K(181)*Y(32)+Y(8)))  
GO TO 100  
39 CONTINUE  
RMOV=(+K(36)*Y(7)+K(37)*Y(13)+K(163)*Y(164)*Y(20))  
FARM=(+K(162)*Y(4)*Y(8)*Y(57))  
GO TO 100  
40 CONTINUE  
RMOV=(+K(50)*Y(18)+K(156)*Y(30))  
FARM=(+K(35)*Y(4)*Y(19))  
GO TO 100  
41 CONTINUE  
RMOV=(+K(24)*Y(30)+K(62)*Y(35)+K(63)*Y(36)+K(84)*Y(37)+K(170)*Y(13)  
1)*Y(57)+K(178)*Y(23)*Y(23))  
FARM=(+K(54)*Y(50)*Y(13))  
GO TO 100  
42 CONTINUE  
RMOV=(Y(57)*(+K(51)*Y(25)+K(172)*Y(13))+K(25)*Y(30)+K(65)*Y(32)+K(  
166)*Y(35)+K(67)*Y(36)+K(68)*Y(37)+K(69)*Y(38))  
FARM=(Y(13)*(+K(60)*Y(48)+K(61)*Y(47)+K(170)*Y(41)*Y(57))+K(174)*Y  
1(43)*Y(57))  
GO TO 100  
43 CONTINUE  
RMOV=(Y(57)*(+K(173)*Y(13)+K(174))+K(26)*Y(30)+K(35)+K(71)*Y  
1(36)+K(172)*Y(37)+K(73)*Y(36))  
FARM=(Y(13)*(+K(55)*Y(53)+K(59)*Y(49)+K(172)*Y(42)*Y(57))+K(175)*Y  
GIONEUT 1701  
GIONEUT 1702  
GIONEUT 1703  
GIONEUT 1704  
GIONEUT 1705  
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GIONEUT 1707  
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GIONEUT 1750  
GIONEUT 1751  
GIONEUT 1752  
GIONEUT 1753  
GIONEUT 1754  
GIONEUT 1755  
GIONEUT 1756

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1 (44)*Y(57))
GIONEUT 1757
GO TO 100
GIONEUT 1758
44 CONTINUE
GIONEUT 1759
RMOV=(Y(57)*(+K(52)*Y(13)+K(175))+K(27)*Y(30)+K(74)*Y(35)+K(75)*Y(
GIONEUT 1760
136)+K(76)*Y(37)+K(77)*Y(38))
GIONEUT 1761
FARM=(Y(57)*(+K(173)*Y(43)*Y(13)+K(176)*Y(45)))
GIONEUT 1762
GO TO 100
GIONEUT 1763
45 CONTINUE
GIONEUT 1764
RMOV=(+K(28)*Y(30)+K(78)*Y(35)+K(79)*Y(36)+K(80)*Y(37)+K(81)*Y(38)
1+K(176)*Y(57))
GIONEUT 1765
FARM=(+K(52)*Y(44)*Y(13)*Y(57))
GIONEUT 1766
GO TO 100
GIONEUT 1767
46 CONTINUE
GIONEUT 1768
RMOV=(+K(57)*Y(25))
GIONEUT 1769
FARM=(+K(171)*Y(41)*Y(23)*Y(23))
GIONEUT 1770
GO TO 100
GIONEUT 1771
47 CONTINUE
GIONEUT 1772
RMOV=(+K(29)*Y(30)+K(61)*Y(13))
GIONEUT 1773
FARM=(+K(53)*Y(50)*Y(13))
GIONEUT 1774
GO TO 100
GIONEUT 1775
48 CONTINUE
GIONEUT 1776
RMOV=(Y(13)*(+K(58)*Y(23)+K(60))+K(91)*Y(36)+K(92)*Y(37)+K(93)*Y(3
18))
GIONEUT 1777
FARM=(+K(57)*Y(46)*Y(25))
GIONEUT 1778
GO TO 100
GIONEUT 1779
49 CONTINUE
GIONEUT 1780
RMOV=(+K(30)*Y(30)+K(59)*Y(13)+K(94)*Y(37)+K(95)*Y(38))
GIONEUT 1781
FARM=(+K(51)*Y(42)*Y(25)*Y(57)+K(58)*Y(48)*Y(13)*Y(23))
GIONEUT 1782
GO TO 100
GIONEUT 1783
50 CONTINUE
GIONEUT 1784
RMOV=(+K(53)*Y(13)+K(54)*Y(13))
GIONEUT 1785
FARM=(+K(37)*Y(39)*Y(13))
GIONEUT 1786
GO TO 100
GIONEUT 1787
51 CONTINUE
GIONEUT 1788
RMOV=(+K(82)*Y(35)+K(83)*Y(36)+K(84)*Y(37)+K(157)*Y(30)+K(177)*Y(
13)*Y(57))
GIONEUT 1789
FARM=(Y(13)*(+K(56)*Y(54)+K(166)*Y(57)))
GIONEUT 1790
GO TO 100
GIONEUT 1791
52 CONTINUE
GIONEUT 1792
RMOV=(+K(85)*Y(85)*Y(86)*Y(87)+K(84)*Y(83)+K(158)*Y(30)+K(178)*Y(
13)*Y(57))
GIONEUT 1793
FARM=(+K(177)*Y(51)*Y(13)*Y(57))
GIONEUT 1794
GO TO 100
GIONEUT 1795
53 CONTINUE
GIONEUT 1796
RMOV=(+K(55)*Y(13)+K(88)*Y(35)+K(89)*Y(36)+K(90)*Y(37)+K(159)*Y(
1))
GIONEUT 1797
FARM=(+K(178)*Y(52)*Y(13)*Y(57))
GIONEUT 1798
GO TO 100
GIONEUT 1799
54 CONTINUE
GIONEUT 1800
RMOV=(+K(56)*Y(13))
GIONEUT 1801
FARM=(+K(165)*Y(5)*Y(25)*Y(23))
GIONEUT 1802
GO TO 100
GIONEUT 1803
55 CONTINUE
GIONEUT 1804
RMOV=(+K(20)*Y(7)+K(21)*Y(10)+K(202))
GIONEUT 1805
FARM=(+K(55)*Y(53)*Y(13)+K(91)*Y(48)*Y(36))
GIONEUT 1806
GO TO 100
GIONEUT 1807
56 CONTINUE
GIONEUT 1808
RMOV=(+K(20)*Y(7)+K(21)*Y(10)+K(202))
GIONEUT 1809
FARM=(+K(55)*Y(53)*Y(13)+K(91)*Y(48)*Y(36))
GIONEUT 1810
GO TO 100
GIONEUT 1811
56 CONTINUE
GIONEUT 1812
GIONEUT 1813

```

400 RMOV=(+K(122)*Y(18)+K(205))
 FARM=(+K(123)*Y(36))
100 FORM(IALT)=FARM
 REMV(IALT)=RMOV
 RETURN
 END
405

GIONEUT 1814
GIONEUT 1815
GIONEUT 1816
GIONEUT 1817
GIONEUT 1818
GIONEUT 1819

```

1      SUBROUTINE RATECN(IALT)
C THIS SUBROUTINE COMPUTES THE IONIZATION RATES. ON THE INITIAL
C ENTRY IT COMPUTES AND SAVES HoF AT NOON FOR O2, N2, AND O.
5      C
C DIMENSION SX(6),QE(6),EX(6),PHI(16),SIGI(16),SIGEFF(16)
C DIMENSION DISOC(:4),XIN(8),ZS(8),CHIS(4)
C COMMON N(692,30),CONC(424,26),U(692,26),NZERO(692),UZERO(692)
10     C
C COMMON FORM(692),REMV(692),E(692),F(692),Z(692),
2      DEDY(692),TEMP(692),DTENDZ(692),DELTAZ(692),SUNSET(692)
C
C COMMON GRAV,COSD,SIND,TIME,CXI,RADIUS,ISPECI,KSPECI,NREAC,K,
1      K2,KMIN1,IITURB,CXINOON,SCXI,NREAC2
C
C COMMON/IONS/WAVE(88),FLUX(88),SIGMA(9,88),R(B,4,8)
C COMMON/CONSTAN/PI,RADSEC,DEGRAD,RADSEC,BOLTZ,SIN21,ATCON
20     C
C COMMON/RATECON/AK(88),BK(88),CK(88),DK(215),CYMB(6,215)
C COMMON/ATCONS/ATWT(26),MASS(26),THERM(26),SYMBOL(56),POLAR(26)
C
C COMMON/LOGIC/NIGHT,DAY,JLOOK,RESET,END
25     C
C COMMON/COLUMN/02COL,03COL,DIS(19,59)
C
C REAL N2NOON,N,MASS,LYALFA,LYBETA
30     C
C COMMON/ALTITUD/ZBOT,ZTOP,TURB,ZTRANS
C
C LOGICAL NIGHT,DAY,JLOOK
35     C
C THE FOLLOWING THREE DATA STATEMENTS CONTAIN THE SOLAR FLUX (X10-9)
C THE IONIZATION CROSS SECTIONS AND THE EFFECTIVE ABSORPTION CROSS
C SECTIONS (X10+18) FOR O2(1DG).
C
40     DATA PHI/0.123,0.202,0.18,0.06,0.16,0.208,0.432,0.092,0.626,0.59,
1      0.08,0.08,0.11,1.63,2.29,0.13/
C
45     DATA SIGI/2.0,3.0,4.0,3.6,4.0,10.0,4.0,4.0,5.0,3.0,5.0,3.8,
1      5.5,5.0,5.0/
C
45     DATA SIGEFF/0.186,0.04,0.021,1.0,0.1,1.0,0.12,1.0,0.2,0.0,4,
1      0.7,0.2,1.15,1.12,1.0/
C
50     C
C THE FOLLOWING TWO DATA STATEMENTS CONTAIN THE ALTITUDES AND THE
C SOLAR ZENITH ANGLE ARRAYS CORRESPONDING TO THE INPUT ARRAYS OF
C THE RATE COEFFICIENTS FOR THE ENERGETIC ELECTRONS.
C
50     DATA ZS/95.,100.,130.,170.,210.,250.,300.,400./
C
55     DATA CHIS/49.77,61.50,76.22,109.7976/
C
55     DATA COSMIC/1.5E-7/
C
55     DATA IPASS/1/

```

```

C      GRAVCOR(ZZ)=1.0/((1.0+ZZ/RADIUS)**2)
60      C      T1=TEMP(IALT)
          ZZ=Z(IALT)
          T2=T1/360.
          D2CBL=0.
65      C      O3COL=0.
          DO 1 J=1,NREAC2
          A1=AK(J)
          B1=BK(J)
          C1=CK(J)
          IF(B1 .NE. 0.) A1=A1*(T2**B1)
          IF(C1 .NE. 0.) A1=A1*EXP(C1/T1)
          DK(J+NREAC)=A1
70      C      CONTINUE
          IF(IPASS .NE. 1) GO TO 30
          DO 26 L=1,16
          PHI(L)=PHI(L)*1.0E+09
          SIGI(L)=SIGI(L)*1.0E-18
          SIGEFF(L)=SIGEFF(L)*1.0E-18
          SIGEFF(L)=SIGEFF(L)*1.0E-18
          26      C      CONTINUE
          IPASS=2
          DO 5 L=1,8
          DO 5 I=1,4
          DO 5 J=1,8
          IF(R(J,I,L) .EQ. 0.) GO TO 3
          R(J,I,L)=ALOG(R(J,I,L))
          GO TO 5
          R(J,I,L)=300.
          3      C      CONTINUE
          5      C      CONTINUE
          C      SAVE H+F AT NOONTIME FOR Q2, O, AND N2.
80      C      GRAV=GRAV*GRAVCOR(ZZ)
          T=BOLTZ*TEMP(IALT)/GRAVITY
          DO 20 I=1,3
          20      C      CALL CHAPMAN(CXNNDN,ZZ,RADIUS,SCALE,APPROX)
          PR=SCALE*APPROX
          IF(I .EQ. 1) Q2NNDN=PR
          IF(I .EQ. 2) N2NNDN=PR
          IF(I .EQ. 3) Q3NNDN=PR
          20      C      CONTINUE
          IF(NIGHT) GO TO 95
          IF(DAY ) GO TO 10
          IF(ZZ .LT. ZTRANS) GO TO 11
          C      GO GET PHOTODISSOCIATION RATE COEFFICIENTS.
105     C      CALL DISSOC(DISOC,IALT)
          DK(192)=DISOC(2)
          DK(193)=DISOC(1)
          DK(194)=DISOC(6)
110     C      GIONEUT 1877
          GIONEUT 1878
          GIONEUT 1879
          GIONEUT 1880
          GIONEUT 1881
          GIONEUT 1882
          GIONEUT 1883
          GIONEUT 1884
          GIONEUT 1885
          GIONEUT 1886
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          GIONEUT 1894
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          GIONEUT 1901
          GIONEUT 1902
          GIONEUT 1903
          GIONEUT 1904
          GIONEUT 1905
          GIONEUT 1906
          GIONEUT 1907
          GIONEUT 1908
          GIONEUT 1909
          GIONEUT 1910
          GIONEUT 1911
          GIONEUT 1912
          GIONEUT 1913
          GIONEUT 1914
          GIONEUT 1915
          GIONEUT 1916
          GIONEUT 1917
          GIONEUT 1918
          GIONEUT 1919
          GIONEUT 1920
          GIONEUT 1921
          GIONEUT 1922
          GIONEUT 1923
          GIONEUT 1924
          GIONEUT 1925
          GIONEUT 1926
          GIONEUT 1927
          GIONEUT 1928
          GIONEUT 1929
          GIONEUT 1930
          GIONEUT 1931
          GIONEUT 1932
          GIONEUT 1933

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115      DK(195)=DISOC(5)
        DK(196)=DISOC(3)
        DK(197)=DISOC(4)
        DK(198)=DISOC(12)
        DK(199)=DISOC(13)
        DK(200)=DISOC(7)
        DK(201)=DISOC(8)
        DK(202)=DISOC(9)
        DK(203)=DISOC(10)
        DK(204)=DISOC(11)
        DK(205)=DISOC(14)

120      IF(.NOT. JLOOK) GO TO 52
        IF(IALT .EQ. 1) GO TO 35
        IF(IALT .EQ. K) GO TO 40
        IF(MOD(IALT-1,12) .NE. 0) GO TO 52
        IA=IALT/12+1
        GO TO 45

125      11     IA=IALT/12+2
        40
        45     DO 50 J=1,14
        DIS(J,IA)=DK(J+191)
        CONTINUE
        DIS(15,IA)=Z(IALT)
        DIS(16,IA)=N(IALT,8)
        DIS(17,IA)=O2COL
        DIS(18,IA)=N(IALT,9)
        DIS(19,IA)=O3COL
        IF(22 .LT. ZTRANS) GO TO 95
        C   COMPUTE RATE COEFFICIENTS FOR ENERGETIC ELECTRON REACTIONS.
        C
130      35
135      50
140      52
145      IF(ZZ .LT. ZS(1)) GO TO 280
        DO 200 I=1,4
        IF(SCXI .EQ. CHIS(I)) GO TO 240
        IF(SCXI .LT. CHIS(I)) GO TO 210
        CONTINUE
        DO 220 J=1,8
        IF(ZZ .EQ. ZS(J)) GO TO 270
        CONTINUE
        DO 235 L=1,8
        DO 225 J=1,8
        XIN(J)=ZS(J)
        YIN(J)=R(J,I-1,L)
        CONTINUE
        A1=YINT(ZZ,XIN,YIN,0)
        B1=YINT(ZZ,XIN,YIN,8)
        DO 230 J=1,8
        YIN(J)=R(J,I,L)
        CONTINUE
        A1=YINT(ZZ,XIN,YIN,0)
        C1=YINT(ZZ,XIN,YIN,8)
        A1=(SCXI-CHIS(I-1))/(CHIS(I)-CHIS(I-1))
        DK(L+207)=EXP(B1-A1*(B1-C1))
        CONTINUE
        DO 1055
        240     DO 245 J=1,8
        IF(ZZ .EQ. ZS(J)) GO TO 260
        CONTINUE
        245

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GIONEUT 1934
 GIONEUT 1935
 GIONEUT 1936
 GIONEUT 1937
 GIONEUT 1938
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 GIONEUT 1940
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 GIONEUT 1981
 GIONEUT 1982
 GIONEUT 1983
 GIONEUT 1984
 GIONEUT 1985
 GIONEUT 1986
 GIONEUT 1987
 GIONEUT 1988
 GIONEUT 1989
 GIONEUT 1990

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DO 255 L=1,8
DO 250 J=1,8
XIN(J)=ZS(J)
YIN(J)=R(J,I,L)
175      CONTINUE
250      AI=YINT(ZZ,XIN,YIN,0)
                  DK(L+207)=EXP(YINT(ZZ,XIN,YIN,0))
255      CONTINUE
180      GO TO 55
260      DO 265 L=1,8
                  DK(L+207)=EXP(R(J,I,L))
265      CONTINUE
185      GO TO 55
270      DO 275 L=1,8
                  A1=(SCXI-CHIS(I-1))/(CHIS(I)-CHIS(I-1))
                  DK(L+207)=EXP(R(J,I-1,L)-A1*(R(J,I-1,L)-R(J,I,L)))
275      CONTINUE
190      GO TO 55
280      DO 285 L=1,8
                  DK(L+207)=0
285      CONTINUE
C      COMPUTE N*H*F FOR O2, N2, AND O AT SOLAR ZENITH ANGLE CXI.
195      C
55      GRAVITY=GRAV*GRAVCOR(ZZ)
                  T=BOLTZ*TEMP(IALT)/GRAVITY
                  DO 60 I=1,3
                  SCALE=T/F(I)
                  CALL CHAPMAN(CXI,ZZ,RADIUS,SCALE,APPROX)
200      IF(I .EQ. 1) E(I)=N(IALT,B)
                  IF(I .EQ. 2) E(I)=N(IALT,C)
                  IF(I .EQ. 3) E(I)=N(IALT,D)
                  EX(I)=E(I)*SCALE*APPROX
205      CONTINUE
60      DO 65 I=1,6
                  QE(I)=0.
                  SX(I)=0.
210      CONTINUE
215      DO 90 L=1,88
                  ABS=0.
                  DO 70 I=1,3
                  ABS=ABS+SIGMA(I,L)*EX(I)
                  CONTINUE
                  AFLUX=0.
                  IF(AFLUX .GT. 650.) GO TO 71
                  AFLUX=FLUX(L)*EXP(-ABS)
220      CONTINUE
225      C      COMPUTE THE IONIZATION COEFFICIENTS. SKIP LYMAN ALPHA LINE.
C
71      IF(L .EQ. 1) GO TO 86
                  IF(L .GE. 74) C0 = C
                  DO 75 I=1,6
                  QE(I)=C5(I)+SIGMA(I+3,L)*AFLUX

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      75    CONTINUE
      80    GO TO 90
      80    DO 85 I=1,6
      85    SX(I)=SX(I)+SIGMA(I+3,L)*AFLUX
      85    CONTINUE
      85    GO TO 90
      86    AFLUX=SFLUX
      86    CONTINUE
      C   PARTITION THE X-RAY IONIZATION ACCORDING TO SWIDER, REVIEWS OF
      C   GEOPHYSICS, 7, PG 573, 1969.
      C
      C   AX=N(IALT,8)+0.5*N(IALT,7)
      C   BX=1.0/(1.0+(N(IALT,23)/(1.15*AX)))
      C   CX=0.71*N(IALT,9)*AX
      C   XTOTAL=SX(1)*N(IALT,8)+SX(2)*N(IALT,23)+SX(3)*N(IALT,7)
      240   1          +SX(4)*N(IALT,11)+SX(5)*N(IALT,21)+SX(6)*N(IALT,16)
      C
      C   COMPUTE TOTAL EUV AND X-RAY IONIZATION RATE COEFFICIENT FOR EACH ION.
      C
      C   DK(185)=QE(3)+SX(3)/N(IALT,7)
      C   DK(186)=QE(1)+SX(1)/N(IALT,8)
      C   DK(187)=QE(2)+SX(2)/N(IALT,23)
      C   DK(189)=QE(6)+SX(6)/N(IALT,16)
      C   DK(190)=QE(4)+SX(4)
      C   DK(191)=QE(5)+SX(5)
      245   255
      C   COMPUTE IONIZATION RATE COEFFICIENT FOR NO BY LYMAN ALPHA.
      C
      C   DK(188)=2.08E-18*SFLUX
      C
      C   COMPUTE IONIZATION RATE COEFFICIENT FOR O2(1DG) FOR WAVELENGTHS
      C   1027-1117 A.
      C
      C   DO 92 L=1,16
      C   SFLUX=SIGEFF(L)*EX(1)
      C   IF(SFLUX .GT. 650.) GO TO 92
      C   DK(207)=DK(207)+PHI(L)*SIG(L)*EXP(-SFLUX)
      C   CONTINUE
      C
      C   ADD CONTRIBUTIONS FROM SCATTERED RADIATIONS AT LYMAN ALPHA.
      C   LYMAN BETA, HE(I)(5B4A), AND HE(II)(304A).
      C
      C   AA=N(IALT,8)*C2NOON
      C   AB=N(IALT,23)*N2NOON
      C   AC=N(IALT,7)*DNODDN
      C   POWER=AA*SIGMAL1,1,1
      C   EXPON=C
      250   255
      C   IF(POWER .GT. 650.) GO TO 91
      C   EXPON=EXP(-POWER)
      C   LYALPH1=1.0E-0212*C2NOON*6.6FLUX(1)*EXPON
      C   POWER=AA*SIGMA(1,2)+A5*SFLUX(1,2)*EXPON
      C   EXPON=C

```

R(1)=K(1)*Y(7)*Y(17)	GIONEUT	2595
R(2)=K(2)*Y(7)*Y(10)	GIONEUT	2596
R(3)=K(3)*Y(7)*Y(12)	GIONEUT	2597
R(4)=K(4)*Y(7)*Y(19)	GIONEUT	2598
R(5)=K(5)*Y(11)*Y(12)	GIONEUT	2599
R(6)=K(6)*Y(11)*Y(12)	GIONEUT	2600
R(7)=K(7)*Y(11)*Y(9)	GIONEUT	2601
R(8)=K(8)*Y(16)*Y(9)	GIONEUT	2602
R(9)=K(9)*Y(16)*Y(19)	GIONEUT	2603
R(10)=K(10)*Y(16)*Y(19)	GIONEUT	2604
R(11)=K(11)*Y(16)*Y(19)	GIONEUT	2605
R(12)=K(12)*Y(10)*Y(10)	GIONEUT	2606
R(13)=K(13)*Y(10)*Y(12)	GIONEUT	2607
R(14)=K(14)*Y(17)*Y(8)	GIONEUT	2608
R(15)=K(15)*Y(27)*Y(13)	GIONEUT	2609
R(16)=K(16)*Y(27)*Y(57)	GIONEUT	2610
R(17)=K(17)*Y(27)*Y(15)	GIONEUT	2611
R(18)=K(18)*Y(20)*Y(57)	GIONEUT	2612
R(19)=K(19)*Y(20)	GIONEUT	2613
R(20)=K(20)*Y(55)*Y(7)	GIONEUT	2614
R(21)=K(21)*Y(55)*Y(10)	GIONEUT	2615
R(22)=K(22)*Y(28)*Y(30)	GIONEUT	2616
R(23)=K(23)*Y(28)*Y(30)	GIONEUT	2617
R(24)=K(24)*Y(41)*Y(30)	GIONEUT	2618
R(25)=K(25)*Y(42)*Y(30)	GIONEUT	2619
R(26)=K(26)*Y(43)*Y(30)	GIONEUT	2620
R(27)=K(27)*Y(44)*Y(30)	GIONEUT	2621
R(28)=K(28)*Y(45)*Y(30)	GIONEUT	2622
R(29)=K(29)*Y(47)*Y(30)	GIONEUT	2623
R(30)=K(30)*Y(49)*Y(30)	GIONEUT	2624
R(31)=K(31)*Y(7)*Y(30)	GIONEUT	2625
R(32)=K(32)*Y(8)*Y(30)	GIONEUT	2626
R(33)=K(33)*Y(4)*Y(18)	GIONEUT	2627
R(34)=K(34)*Y(4)*Y(16)	GIONEUT	2628
R(35)=K(35)*Y(4)*Y(19)	GIONEUT	2629
R(36)=K(36)*Y(39)*Y(7)	GIONEUT	2630
R(37)=K(37)*Y(39)*Y(13)	GIONEUT	2631
R(38)=K(38)*Y(28)*Y(7)	GIONEUT	2632
R(39)=K(39)*Y(28)*Y(7)	GIONEUT	2633
R(40)=K(40)*Y(3)*Y(11)	GIONEUT	2634
R(41)=K(41)*Y(1)*Y(7)	GIONEUT	2635
R(42)=K(42)*Y(6)*Y(8)	GIONEUT	2636
R(43)=K(43)*Y(6)*Y(8)	GIONEUT	2637
R(44)=K(44)*Y(6)*Y(8)	GIONEUT	2638
R(45)=K(45)*Y(2)*Y(8)	GIONEUT	2639
R(46)=K(46)*Y(2)*Y(8)	GIONEUT	2640
R(47)=K(47)*Y(2)*Y(23)	GIONEUT	2641
R(48)=K(48)*Y(2)*Y(23)	GIONEUT	2642
R(49)=K(49)*Y(2)*Y(18)	GIONEUT	2643
R(50)=K(50)*Y(40)*Y(18)	GIONEUT	2644
R(51)=K(51)*Y(42)*Y(25)*Y(57)	GIONEUT	2645
R(52)=K(52)*Y(44)*Y(13)*Y(57)	GIONEUT	2646
R(53)=K(53)*Y(50)*Y(13)	GIONEUT	2647
R(54)=K(54)*Y(50)*Y(13)	GIONEUT	2648
R(55)=K(55)*Y(53)*Y(13)	GIONEUT	2649
R(56)=K(56)*Y(54)*Y(13)	GIONEUT	2650
R(57)=K(57)*Y(46)*Y(25)	GIONEUT	2651

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1      SUBROUTINE CHEMPR(IZ)
2
3      C THIS SUBROUTINE PRINTS OUT THE FULL CHEMISTRY LIST ALONG WITH
4      C THE RATE COEFFICIENTS AND THE RATES OF ALL REACTIONS AT THE
5      C ALTITUDE CORRESPONDING TO THE INDEX IZ.
6
7      C DIMENSION Y(57),R(215),K(215)
8
9      C COMMON N(692,30),CONC(424,26),U(692,26),NZERO(692),UZERO(692)
10     C COMMON FORM(692),RENV(692),E(692),F(692),Z(692),
11        2       DEDY(692),TEMP(692),DTENDZ(692),DELTAZ(692),SUNSET(692)
12
13     C COMMON GRAV,COSD,SIND,TIME,CXI,RADIUS,ISPECI,KSPECI,NREAC,KTOP,
14        1       K2,KMIN1,ITURB,CXIND0N,SCXI,NREAC2
15
16     C COMMON/ESM/OLDN(692,30),OLDCON(424,26),OLDU(692,26),RATES(88,692)
17
18     C COMMON/LOGIC/NIGHT,DAY,JLDOK,RESET,END
19
20     C COMMON/RATECON/AK(88),BK(88),CK(88),DK(215),CYMB(6,215)
21
22     C LEVEL 3, OLDN,OLDCON,OLDU,RATES
23     C REAL K,N,NZERO
24     C LOGICAL NIGHT
25     C EQUIVALENCE (Y,E),(R,F),(NZERO,K)
26
27     C WRITE(6,1) Z(IZ),SCXI
28     1   FORMAT(1H1, * CHEMISTRY AT *,F7.2, * KM. AND AT *, 1PE12.5)
29
30     C WRITE(6,5)
31     5   FORMAT(1H0,1BX,*CHEMICAL REACTION*,.35X,*RATE CONSTANT*,.5X,
32           1*FORWARD RATE/(CM3/SEC*)*)
33     C DO 6 I=1,NREAC
34     C   K(I)=DK(I)
35
36     C CONTINUE
37     C IF(.NOT. NIGHT) GO TO 10
38     C   K(96)=K(99)=K(104)=K(105)=K(110)=0.
39     C   K(113)=K(114)=K(118)=K(120)=K(123)=0.
40     C   CALL MOLEV(RATES(1,IZ),K(128),88)
41     C   IF(IZ .LE. K2) GO TO 11
42     C   K(31)=K(32)=K(35)=0.
43     C   K(160)=K(161)=K(162)=K(165)=K(166)=0.
44
45     C   DO 25 I=1,ISPECI
46     C     IF(I .GT. 30) GO TO 4
47     C     Y(I)=N(IZ,I)
48     C     GO TO 25
49
50     C   4   IF(IZ .GT. K2) GO TO 15
51     C     Y(I)=CONC(IZ,I-30)
52     C     GO TO 25
53
54     C   15   Y(I)=0.
55     C   25   CONTINUE
56     C     Y(57)=0.
57     C     DO 30 I=1,ISPECI
58     C       Y(57)=Y(57)+Y(I)
59     C     CONTINUE
60
61     C   30

```

```

1      SUBROUTINE INDEX(COL,LIM,I,DEL)
C   THIS SUBROUTINE COMPUTES AN INDEX I AND AN INCREMENT DEL FROM
C   THE COLUMN DENSITY COL.
5      C
      REAL LOG2,LOG5
      LOGICAL ISKIP
      DATA ISKIP/.FALSE./
      C
      IF(ISKIP) GO TO 5
      LOG2= ALOG10(2.0)
      LOG5= ALOG10(5.0)
      ISKIP=.TRUE.
      C
      A=ALOG10(COL)
      INTA=INT(A)
      FA=A-INTA
      I=3*(INTA-LIM)+1
      IF(FA.GT.LOG2) GO TO 10
      RETURN
      DEL=FA/LOG2
      10     IF(FA.GT. LOG5) GO TO 20
      DEL=(FA-LOG2)/(LOG5-LOG2)
      I=I+1
      RETURN
      20     DEL=(FA-LOG5)/(1.0-LOG5)
      I=I+2
      RETURN
      END
      GIONEUT 2509
      GIONEUT 2510
      GIONEUT 2511
      GIONEUT 2512
      GIONEUT 2513
      GIONEUT 2514
      GIONEUT 2515
      GIONEUT 2516
      GIONEUT 2517
      GIONEUT 2518
      GIONEUT 2519
      GIONEUT 2520
      GIONEUT 2521
      GIONEUT 2522
      GIONEUT 2523
      GIONEUT 2524
      GIONEUT 2525
      GIONEUT 2526
      GIONEUT 2527
      GIONEUT 2528
      GIONEUT 2529
      GIONEUT 2530
      GIONEUT 2531
      GIONEUT 2532
      GIONEUT 2533
      GIONEUT 2534
      GIONEUT 2535
      GIONEUT 2536
      GIONEUT 2537

```

```

1      FUNCTION TRPLT1(ARRAY,103,DELO3)
C   THIS FUNCTION INTERPOLATES IN THE ONE DIMENSIONAL PHOTODISSOCIATION
C   RATE COEFFICIENT TABLES.
C
5      C   DIMENSION ARRAY(1)
C
C   A=ARRAY(103)
C   B=ARRAY(103+1)
C   TRPLT1=10.0***(A-DELO3*(A-B))
C   RETURN
10     END

```

GIONEUT 2497
 GIONEUT 2498
 GIONEUT 2499
 GIONEUT 2500
 GIONEUT 2501
 GIONEUT 2502
 GIONEUT 2503
 GIONEUT 2504
 GIONEUT 2505
 GIONEUT 2506
 GIONEUT 2507
 GIONEUT 2508

```

1      FUNCTION TRPLT2(ARRAY,N,M,I02,DEL02,I03,DEL03)
2
3      C THIS FUNCTION INTERPOLATES IN THE TWO DIMENSIONAL PHOTODISSOCIATION
4      C RATE COEFFICIENT TABLES.
5      C
6      C DIMENSION ARRAY(N,M)
7
8      C
9      A=ARRAY(I02,I03)
10     B=ARRAY(I02,I03+1)
11     C=ARRAY(I02+1,I03)
12     D=ARRAY(I02+1,I03+1)
13     Y1=A-DEL03*(A-B)
14     Y2=C-DEL03*(C-D)
15     TRPLT2=10.0**((Y1-DEL02)*(Y1-Y2))
16     RETURN
17     END

```

230 IF(02COL .GT. 23.5) GO TO 210
 DO 200 II=1,12
 IF(02COL .LT. 02C(II)) GO TO 215
 CONTINUE
 STOP 7
200 205 DISOC(13)=10.0** (PNO(1))
 GO TO 220
210 DISOC(13)=0.
 GO TO 220
215 DEL02=(02COL-02C(II-1))/(02C(II)-02C(II-1))
 DISOC(13)=10.0** (PNO(II-1)-DEL02*(PNO(II-1)-PNO(II)))
220 02COL=10.0** (02COL)
 RETURN
 END
240 245 GIONEUT 2467
 GIONEUT 2468
 GIONEUT 2469
 GIONEUT 2470
 GIONEUT 2471
 GIONEUT 2472
 GIONEUT 2473
 GIONEUT 2474
 GIONEUT 2475
 GIONEUT 2476
 GIONEUT 2477
 GIONEUT 2478
 GIONEUT 2479
 GIONEUT 2480

```

175      IF(03COLM .GT. 1.0E+21) 03COL=1.0E+21
         LIMIT=14
         CALL INDEX(03COL,LIMIT,103,DEL03)
         DISOC(9)=TRPLT1(J202,103,DEL03)

C   REACTION CO2 + HV = CO + O1D IN DISOC(10)
C
C   Q2COL=02COLM
C   1F(02COLM .LT. 1.0E+13) 02COL=1.1E+13
C   1F(02COLM .GT. 5.0E+22) 02COL=5.0E+22
         LIMIT=13
         CALL INDEX(02COL,LIMIT,102,DEL02)
         Q3COL=03COLM
         1F(03COLM .LT. 1.0E+11) 03COL=1.1E+11
         1F(03COLM .GT. 2.0E+20) 03COL=2.0E+20
         LIMIT=11
         CALL INDEX(03COL,LIMIT,103,DEL03)
         DISOC(10)=TRPLT2(J203,30,29,102,DEL02,103,DEL03)

180      C   REACTION CO2 + HV = CO + O IN DISOC(11)
C
C   Q2COL=02COLM
C   1F(02COLM .LT. 1.0E+13) 02COL=1.1E+13
C   1F(02COLM .GT. 2.0E+24) 02COL=2.0E+24
         LIMIT=13
         CALL INDEX(02COL,LIMIT,102,DEL02)
         Q3COL=03COLM
         1F(03COLM .LT. 1.0E+12) 03COL=1.1E+12
         1F(03COLM .GT. 1.0E+21) 03COL=1.0E+21
         LIMIT=12
         CALL INDEX(03COL,LIMIT,103,DEL03)
         DISOC(11)=TRPLT2(J204,35,28,102,DEL02,103,DEL03)

185      C   REACTION CO2 + HV = CO + O IN DISOC(12)
C
C   Q2COL=02COLM
C   1F(02COLM .LT. 1.0E+13) 02COL=1.1E+13
C   1F(02COLM .GT. 2.0E+24) 02COL=2.0E+24
         LIMIT=13
         CALL INDEX(02COL,LIMIT,102,DEL02)
         Q3COL=03COLM
         1F(03COLM .LT. 1.0E+12) 03COL=1.1E+12
         1F(03COLM .GT. 1.0E+21) 03COL=1.0E+21
         LIMIT=12
         CALL INDEX(03COL,LIMIT,103,DEL03)
         DISOC(12)=TRPLT2(J205,30,28,102,DEL02,103,DEL03)

190      C   REACTION H02 + HV = O2 + H IN DISOC(13)
C
C   Q2COL=02COLM
C   1F(02COLM .LT. 1.0E+17) 02COL=1.1E+17
C   1F(02COLM .GT. 2.0E+25) 02COL=2.0E+25
         LIMIT=17
         CALL INDEX(02COL,LIMIT,102,DEL02)
         Q3COL=03COLM
         1F(03COLM .LT. 1.0E+12) 03COL=1.1E+12
         1F(03COLM .GT. 1.0E+21) 03COL=1.0E+21
         LIMIT=12
         CALL INDEX(03COL,LIMIT,103,DEL03)
         DISOC(13)=TRPLT2(J206,26,28,102,DEL02,103,DEL03)

195      C   REACTION H02 + HV = O2 + H IN DISOC(14)
C
C   Q2COL=02COLM
C   1F(02COLM .LT. 1.0E+17) 02COL=1.1E+17
C   1F(02COLM .GT. 2.0E+25) 02COL=2.0E+25
         LIMIT=17
         CALL INDEX(02COL,LIMIT,102,DEL02)
         Q3COL=03COLM
         1F(03COLM .LT. 1.0E+12) 03COL=1.1E+12
         1F(03COLM .GT. 1.0E+21) 03COL=1.0E+21
         LIMIT=12
         CALL INDEX(03COL,LIMIT,103,DEL03)
         DISOC(14)=TRPLT2(J207,26,28,102,DEL02,103,DEL03)

200      C   REACTION NO3 + HV = NO2 + O2 IN DISOC(15)
C
C   Q2COL=02COLM
C   1F(02COLM .LT. 1.0E+14) 03COL=1.1E+14
C   1F(02COLM .GT. 1.0E+21) 03COL=1.0E+21
         LIMIT=14
         CALL INDEX(03COL,LIMIT,103,DEL03)
         DISOC(15)=TRPLT1(J208,26,28,102,DEL02,103,DEL03)
         Q2COL=ALOG10(02COLM)
         1F(02COLM .LT. 18.0) GO TO 205

```

```

115      C DISOC(4)=TRPLT2(J197,27,28,102,DEL02,103,DEL03)
      C REACTION O3 + HV = O1D + O2ID IN DISOC(5)
      C
      C 02COL=02COLM
      C   IF(02COLM .LT. 1.0E+15) 02COL=1.1E+15
      C   IF(02COLM .GT. 2.0E+25) 02COL=2.0E+25
      C   LIMIT=15
      C   CALL INDEX(02COL,LIMIT,102,DEL02)
      C
      C 03COL=03COLM
      C   IF(03COLM .LT. 1.0E+12) 03COL=1.1E+12
      C   IF(03COLM .GT. 1.0E+21) 03COL=1.0E+21
      C   LIMIT=12
      C   CALL INDEX(03COL,LIMIT,103,DEL03)
      C   DISOC(5)=TRPLT2(J195,32,28,102,DEL02,103,DEL03)

120      C REACTION O3 + HV = O + O2 IN DISOC(6)
      C
      C 03COL=03COLM
      C   IF(03COLM .LT. 1.0E+14) 03COL=1.1E+14
      C   IF(03COLM .GT. 1.0E+21) 03COL=1.0E+21
      C   LIMIT=14
      C   CALL INDEX(03COL,LIMIT,103,DEL03)
      C   DISOC(6)=TRPLT1(J194,103,DEL03)

125      C REACTION NO2 + HV = NO + O2 IN DISOC(7)
      C
      C 02COL=02COLM
      C   IF(02COLM .LT. 1.0E+19) 02COL=1.1E+19
      C   IF(02COLM .GT. 1.0E+25) 02COL=1.0E+25
      C   LIMIT=19
      C   CALL INDEX(02COL,LIMIT,102,DEL02)
      C
      C 03COL=03COLM
      C   IF(03COLM .LT. 1.0E+14) 03COL=1.1E+14
      C   IF(03COLM .GT. 5.0E+18) 03COL=5.0E+18
      C   LIMIT=14
      C   CALL INDEX(03COL,LIMIT,103,DEL03)
      C   DISOC(7)=TRPLT2(J200,19,15,102,DEL02,103,DEL03)

130      C REACTION N2O + HV = N2 + O1D IN DISOC(8)
      C
      C 02COL=02COLM
      C   IF(02COLM .LT. 1.0E+17) 02COL=1.1E+17
      C   IF(02COLM .GT. 2.0E+25) 02COL=2.0E+25
      C   LIMIT=17
      C   CALL INDEX(02COL,LIMIT,102,DEL02)
      C
      C 03COL=03COLM
      C   IF(03COLM .LT. 1.0E+11) 03COL=1.1E+11
      C   IF(03COLM .GT. 5.0E+19) 03COL=5.0E+19
      C   LIMIT=11
      C   CALL INDEX(03COL,LIMIT,103,DEL03)
      C   DISOC(8)=TRPLT2(J201,26,27,102,DEL02,103,DEL03)

135      C REACTION HNO2 + HV = OH + OH IN DISOC(9)
      C
      C 03COL=03COLM
      C   IF(03COLM .LT. 1.0E+14) 03COL=1.1E+14

```

```

1      CONTINUE
35     03COL=03COLM
60     02COL=02COLM
C      REACTION O2 + HV = O + OH IN DISOC(1)
C
C      IF(02COLM .LT. 1.0E+13) 02COL=1.1E+13
C      IF(02COLM .GT. 1.0E+22) 02COL=1.0E+22
C      LIMIT=13
CALL INDEX(02COL,LIMIT,102,DEL02)

C      03COL=03COLM
C      IF(03COLM .LT. 1.0E+12) 03COL=1.1E+12
C      IF(03COLM .GT. 2.0E+19) 03COL=1.9E+19
C      LIMIT=12
CALL INDEX(03COL,LIMIT,103,DEL03)
DISOC(1)=TRPLT2(J193,28,26,102,DEL02,103,DEL03)

75     C      REACTION O2 + HV = O + O IN DISOC(2)
C
C      02COL=02COLM
C      IF(02COLM .LT. 1.0E+17) 02COL=1.1E+17
C      IF(02COLM .GT. 2.0E+25) 02COL=2.0E+25
C      LIMIT=17
CALL INDEX(02COL,LIMIT,102,DEL02)

C      03COL=03COLM
C      IF(03COLM .LT. 1.0E+13) 03COL=1.1E+13
C      IF(03COLM .GT. 1.0E+21) 03COL=1.0E+21
C      LIMIT=13
CALL INDEX(03COL,LIMIT,103,DEL03)
DISOC(2)=TRPLT2(J192,26,25,102,DEL02,103,DEL03)

75     C      REACTION H2O + HV = OH + H IN DISOC(3)
C
C      02COL=C2COLM
C      IF(02COLM .LT. 1.0E+13) 02COL=1.1E+13
C      IF(02COLM .GT. 1.0E+23) 02COL=1.0E+23
C      LIMIT=13
CALL INDEX(02COL,LIMIT,102,DEL02)

C      03COL=03COLM
C      IF(03COLM .LT. 1.0E+11) 03COL=1.1E+11
C      IF(03COLM .GT. 5.0E+20) 03COL=5.0E+20
C      LIMIT=11
CALL INDEX(03COL,LIMIT,103,DEL03)
DISOC(3)=TRPLT2(J196,31,30,102,DEL02,103,DEL03)

90     C      REACTION H2O2 + HV = OH + OH IN DISOC(4)
C
C      02COL=02COLM
C      IF(02COLM .LT. 1.0E+17) 02COL=1.1E+17
C      IF(02COLM .GT. 5.0E+25) 02COL=5.0E+25
C      LIMIT=17
CALL INDEX(02COL,LIMIT,102,DEL02)

C      03COL=C3COLM
C      IF(03COLM .LT. 1.0E+12) 03COL=1.1E+12
C      IF(03COLM .GT. 1.0E+21) 03COL=1.0E+21
C      LIMIT=12
CALL INDEX(C3COL,LIMIT,103,DEL03)

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1      C SUBROUTINE DISSOC(DISDC,IALT)
2      C DIMENSION DISOC(14),O2C(12),PNO(12)
3      C COMMON N(692,30),CONC(424,26),U(692,26),NZERO(692),UZERO(692)
4      C COMMON FORM(692),REMV(692),E(692),F(692),Z(692),
5      C DEDY(692),TEMP(692),DTENDZ(692),DELTAZ(692),SUMSET(692)
6      C COMMON GRAV,COSD,SIND,TIME,CX1,RADIUS,ISPECI,KSPECI,NREAC,X,
7      C K2,KMIN1,ITURB,CXINDON,SCXI,NREAC2
8      C COMMON/ATCONS/AITNT(26),MASS(26),THERM(26),SYMBOL(56),POLAR(26)
9      C COMMON/CONSTAN/PI,RADDEG,DEGRAD,RADSEC,BDLTZ,SIN2I,ATCON
10     C COMMON/JAYS/J192(26,25),J193(28,26),J194(22),J195(32,28),
11     C J196(31,30),J197(27,28),J198(26,28),J200(19,15),
12     C J201(26,27),J202(22),J203(30,29),J204(35,28),J205(22)
13     C COMMON/COLUM/O2CDLM,O3CDLM,DIS(19,59)
14     C REAL N,MASS,J192,J193,J194,J195,J196,J198,J200,J201,J202
15     C REAL J203,J204,J205
16     C LOGICAL JSTART
17     C DATA JSTART/.TRUE./
18
19     C THE FOLLOWING TWO DATA STATEMENTS CONTAIN THE NO PHOTODISSOCIATION
20     C RATE COEFFICIENTS (PNO) AS FUNCTIONS OF THE COLUMN DENSITY OF
21     C O2 (O2C). DATA FROM CIESLIK AND NICOLET, PLANET SPACE SCI.
22     C 21, 925, 1973.
23
24     C DATA O2C/18.0,18.5,19.0,19.5,20.0,20.5,21.0,21.5,22.0,22.5,23.0,
25     C 23.5/
26     C DATA PNO/1.41E-5,1.40E-5,1.38E-5,1.33E-5,1.22E-5,1.02E-5,
27     C 6.98E-6,3.58E-6,8.93E-7,1.26E-7,4.45E-12,1.50E-29/
28
29     C GRAVCOR(ZE)=1.0/((1.0+ZE/RADIUS)**2)
30
31     C COMPUTE COLUMN DENSITIES OF O2 AND O3 USING THE SWIDER
32     C APPROXIMATION TO THE CHAPMAN FUNCTION.
33
34     C ALT=Z(IALT)
35     C T=(BDLTZ*TEMP(IALT))/(GRAV*GRAVCOR(ALT))
36     C SCALE=T/MASS(0)
37     C CALL CHAPMAN(CX1,ALT,RADIUS,SCALE,APPROX)
38     C O2CDLM=N(IALT,B)*SCALE*APPROX
39     C CALL CHAPMAN(CX1,ALT,RADIUS,SCALE,APPROX)
40     C O3CDLM=N(IALT,G)*SCALE*APPROX
41     C IF(.NOT.JSTART) GO TO 35
42     C JSTART=.FALSE.
43
44     C REDUCE THE NO PHOTODISSOCIATION RATE COEFFICIENTS BY FACTOR OF 4.
45
46     DO 1 I=1,12
47     PNO(I)=ALOG10(PNO(I))/4.0

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1          SUBROUTINE CHAPMAN(CXI,Z,R,H,F)
2
3          C THIS SUBROUTINE COMPUTES AN APPROXIMATION FOR THE CHAPMAN
4          C FUNCTION USING EQUATIONS F(2) AND FN OF SNIDER FROM
5          C "ON THE ACCURACY OF CERTAIN APPROXIMATIONS FOR THE CHAPMAN
6          C FUNCTION". AFCRL-67-0468.
7
8          DATA P10V2/1.570796326795/, SQPI2/8.862269254526E-01/
9          DATA PI/3.14159265359/
10         X=(R+Z)/(H*1.0E-05)
11         Y=X*SIN(CXI)
12         IF(CXI.GT.P10V2) GO TO 10
13         W=X-Y
14         IF(W.NE.0.) GO TO 1
15         F=SQRT(1.0+2.0*X)*SQPI2
16         RETURN
17
18         1   A=SQRT(W)
19         IF(W.GE.9.5) GO TO 5
20         COM=ERFC(A)
21         F=-X*COS(CXI)+SQRT(1.0+2.0*X-W)*(A+SQPI2*COM*EXP(W))
22         RETURN
23
24         5   F=-X*COS(CXI)+SQRT(1.0+2.0*X-W)*(A+(1.0/(2.0*A))*(1.0
25           -(1.0/(2.0*W))+(3.0/(4.0*(W**2)))-(15.0/(8.0*(W**3)))))
26
27         10  A=-SQRT(Y/2.0)/TAN(CXI)
28         EF=ERF(A)
29         F=SQRT(P1*Y/2.0)*(1.0+EF)
30         RETURN
31
32         END
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      RETURN
      C   ERROR EXITS.
      C
60
      20   PRINT 902, J,J1,IJ,JH,X(J),X(J1),DX
      902   FORMAT(" ERROR IN YINT.          DX .LE. 0."
1/* J=.14., J1=.14., JH=.14., X(J)=.1PE13.5,
2. X(J1)=.E13.5., DX=.E13.5)
      GO TO 910
      30   PRINT 903, JX,JH,XI,X(JX-1),X(JH)
      903   FORMAT(" ERROR IN YINT.  XI OUTSIDE RANGE OF X."
1/* JX=.14., JH=.14., XI=.1PE13.5., X(JX-1)=.E13.5,
2. X(JH)=.E13.5)
      GO TO 910
      40   PRINT 904, J,J1,IJ,JH,X(J),X(J+1),DX1
      904   FORMAT(" ERROR IN YINT.          DX1 .LE. 0."
1/* J=.14., J1=.14., JH=.14., X(J)=.1PE13.5,
2. X(J+1)=.E13.5., DX1=.E13.5)
      75
      910  STOP
      END

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GIONEUT 2190
GIONEUT 2191
GIONEUT 2192
GIONEUT 2193
GIONEUT 2194
GIONEUT 2195
GIONEUT 2196
GIONEUT 2197
GIONEUT 2198
GIONEUT 2199
GIONEUT 2200
GIONEUT 2201
GIONEUT 2202
GIONEUT 2203
GIONEUT 2204
GIONEUT 2205
GIONEUT 2206
GIONEUT 2207
GIONEUT 2208
GIONEUT 2209

```

1           FUNCTION YINT(XII,X,Y,JHI)
2
3           C THIS FUNCTION RETURNS INTERPOLATED YINT FOR INPUT XII. ARRAY X
4           C MUST BE MONOTONICALLY INCREASING. JHI IS NUMBER OF POINTS IN
5           C X AND Y ARRAYS.
6
7           C INITIALIZE BY CALL WITH JHI=0 BEFORE EACH ARRAY CALL.
8
9           C EXAMPLE
10          C INITIALIZE
11          C YY=YINT(XII,X,Y,0)
12          C LOOP ARRAY INTERPOLATE
13          DO 10 K=1,100
14          XII=K
15          YI(K)=YINT(XII,X,Y,JHI)
16
17          C DIMENSION X(1),Y(1)
18          JH=JHI
19          XI=XII
20          IF(JH .GT. 0) GO TO 8
21
22          C INITIALIZE
23          J1=0
24
25          RETURN
26          IF(J1 .LT. 1) JX=2
27          IF(J1 .LT. 1) GO TO 1
28
29          C TEST IF XII BETWEEN CURRENT X(J1),X(J1+1)
30          IF(XI .GE. X(J1) .AND. XI .LE. X(J1+1)) GO TO 2
31
32          C SEARCH FOR STRADDLING X VALUES
33          DD 10 J=UX,JH
34          J1=J-1
35
36          DX=X(J)-X(J1)
37          C TEST FOR NON-ZERO DX
38          IF(DX .LE. 0.) GO TO 20
39          IF(XI .GE. X(J1) .AND. XI .LE. X(J)) GO TO 4
40          CONTINUE
41          C ERROR
42          GO TO 30
43          C COMPUTE SA=Avg SLOPE BETWEEN X(J1) AND X(J1+1)
44          SA=(Y(J)-Y(J1))/DX
45          JX=J
46
47          C SET PV,NXT SLOPE TO SA TEMPORARILY
48          SAP=SAN=SA
49
50          C SET PVS SLOPE IF J1.GT.1
51          IF(J1 .GT. 1) SAP=(Y(J1)-Y(J1-1))/(X(J1)-X(J1-1))
52          IF(J .GE. JH) GO TO 6
53          DX1=X(J+1)-X(J)
54          IF(DX1 .LE. 0.) GO TO 40
55          C SET NEXT SLOPE IF J1 .LT. JH1-1
56          SAN=(Y(J+1)-Y(J))/DX1
57
58          S1=S2=0.
59          IF(SA*SAP .GT. 0.) S1=2.*SA*SAP/(SA+SAN)
60          IF(SA-SAN .GT. 0.) S2=2.*SA*SAN/(SA+SAN)
61          A2=(3.*SA-2.*S1-S2)/DX
62          A3=(SA-S1-A2*DX)/DX-DX
63
64          C THIS IS CUBIC FIT FOR STRADDLES XI
65          2
66          XI=XI-X(J1)
67          YINT=Y(J1)+DX1*(S1+DX1*(A2+DX1*A3))
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      IF(POWER .GT. 650.) GO TO 96
      EXPON=EXP(-POWER)
      LYBETA=4.0E-03*SIGMA(4.2)*FLUX(2)*EXPON
      POWER=AA*SIGMA(1.55)+AB*SIGMA(2.55)+AC*SIGMA(3.55)
      EXPON=0.
      IF(POWER .GT. 650.) GO TO 100
      EXPON=EXP(-POWER)
      HEI=1.1E-03*FLUX(55)*EXPON
      POWER=AA*SIGMA(1.68)+AB*SIGMA(2.68)+AC*SIGMA(3.68)
      EXPON=0.
      IF(POWER .GT. 650.) GO TO 105
      EXPON=EXP(-POWER)
      HEII=1.0E-03*FLUX(68)*EXPON
      DK(185)=DK(185)+HEI*SIGMA(6.55)+HEII*SIGMA(6.68)
      DK(186)=DK(186)+LYBETA+HEI*SIGMA(4.55)+HEII*SIGMA(4.68)
      DK(187)=DK(187)+HEI*SIGMA(5.55)+HEII*SIGMA(5.68)
      DK(188)=DK(188)+LYALFA
      DK(190)=DK(190)+HEI*SIGMA(7.55)+HEII*SIGMA(7.68)
      DK(191)=DK(191)+HEI*SIGMA(8.55)+HEII*SIGMA(8.68)

290      C   ADD COSMIC RAY IONIZATION.    Q=1.0E-17*N
            C
            DD 110  I=185.191
            DK(I)=DK(I)+COSMIC
            CONTINUE
            DK(206)=DK(185)
            RETURN
            END

300      GIONEUT 2105
            GIONEUT 2106
            GIONEUT 2107
            GIONEUT 2108
            GIONEUT 2109
            GIONEUT 2110
            GIONEUT 2111
            GIONEUT 2112
            GIONEUT 2113
            GIONEUT 2114
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            GIONEUT 2116
            GIONEUT 2117
            GIONEUT 2118
            GIONEUT 2119
            GIONEUT 2120
            GIONEUT 2121
            GIONEUT 2122
            GIONEUT 2123
            GIONEUT 2124
            GIONEUT 2125
            GIONEUT 2126
            GIONEUT 2127
            GIONEUT 2128
            GIONEUT 2129
            GIONEUT 2130
            GIONEUT 2131
            GIONEUT 2132

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115	R(58)=K(58)*Y(48)*Y(13)*Y(23) R(59)=K(59)*Y(49)*Y(13) R(60)=K(60)*Y(48)*Y(13) R(61)=K(61)*Y(47)*Y(13) R(62)=K(62)*Y(41)*Y(35) R(63)=K(63)*Y(41)*Y(36) R(64)=K(64)*Y(41)*Y(37) R(65)=K(65)*Y(42)*Y(32) R(66)=K(66)*Y(42)*Y(35) R(67)=K(67)*Y(42)*Y(36) R(68)=K(68)*Y(42)*Y(37) R(69)=K(69)*Y(42)*Y(38) R(70)=K(70)*Y(43)*Y(35) R(71)=K(71)*Y(43)*Y(36) R(72)=K(72)*Y(43)*Y(37) R(73)=K(73)*Y(43)*Y(38) R(74)=K(74)*Y(44)*Y(35) R(75)=K(75)*Y(44)*Y(36) R(76)=K(76)*Y(44)*Y(37) R(77)=K(77)*Y(44)*Y(38) R(78)=K(78)*Y(45)*Y(35) R(79)=K(79)*Y(45)*Y(36) R(80)=K(80)*Y(45)*Y(37) R(81)=K(81)*Y(45)*Y(38) R(82)=K(82)*Y(51)*Y(35) R(83)=K(83)*Y(51)*Y(36) R(84)=K(84)*Y(51)*Y(37) R(85)=K(85)*Y(52)*Y(35) R(86)=K(86)*Y(52)*Y(36) R(87)=K(87)*Y(52)*Y(37) R(88)=K(88)*Y(53)*Y(35) R(89)=K(89)*Y(53)*Y(36) R(90)=K(90)*Y(53)*Y(37) R(91)=K(91)*Y(48)*Y(36) R(92)=K(92)*Y(48)*Y(37) R(93)=K(93)*Y(48)*Y(38) R(94)=K(94)*Y(49)*Y(37) R(95)=K(95)*Y(49)*Y(38) R(96)=K(96)*Y(31) R(97)=K(97)*Y(31)*Y(7)	GIONEUT 2652 GIONEUT 2653 GIONEUT 2654 GIONEUT 2655 GIONEUT 2656 GIONEUT 2657 GIONEUT 2658 GIONEUT 2659 GIONEUT 2660 GIONEUT 2661 GIONEUT 2662 GIONEUT 2663 GIONEUT 2664 GIONEUT 2665 GIONEUT 2666 GIONEUT 2667 GIONEUT 2668 GIONEUT 2669 GIONEUT 2670 GIONEUT 2671 GIONEUT 2672 GIONEUT 2673 GIONEUT 2674 GIONEUT 2675 GIONEUT 2676 GIONEUT 2677 GIONEUT 2678 GIONEUT 2679 GIONEUT 2680 GIONEUT 2681 GIONEUT 2682 GIONEUT 2683 GIONEUT 2684 GIONEUT 2685 GIONEUT 2686 GIONEUT 2687 GIONEUT 2688 GIONEUT 2689 GIONEUT 2690 GIONEUT 2691 GIONEUT 2692 GIONEUT 2693 GIONEUT 2694 GIONEUT 2695 GIONEUT 2696 GIONEUT 2697 GIONEUT 2698 GIONEUT 2699 GIONEUT 2700 GIONEUT 2701 GIONEUT 2702 GIONEUT 2703 GIONEUT 2704 GIONEUT 2705 GIONEUT 2706 GIONEUT 2707 GIONEUT 2708
120	R(98)=K(98)*Y(31)*Y(9) R(99)=K(99)*Y(32)	
125	R(100)=K(100)*Y(32)*Y(7) R(101)=K(101)*Y(32)*Y(7) R(102)=K(102)*Y(32)*Y(9) R(103)=K(103)*Y(32)*Y(20) R(104)=K(104)*Y(33)	
130	R(105)=K(105)*Y(33) R(106)=K(106)*Y(33)*Y(7) R(107)=K(107)*Y(34)*Y(7)	
135	R(108)=K(108)*Y(34)*Y(7) R(109)=K(109)*Y(34)*Y(25)	
140	R(110)=K(110)*Y(37) R(111)=K(111)*Y(37)*Y(7)	
145	R(112)=K(112)*Y(37)*Y(19) R(113)=K(113)*Y(38) R(114)=K(114)*Y(38)	
150		
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			R((115)=K(115)*Y(38)*Y(7)	GIONEUT	2709
			R((116)=K(116)*Y(38)*Y(9)	GIONEUT	2710
			R((117)=K(117)*Y(38)*Y(18)	GIONEUT	2711
			R((118)=K(118)*Y(35)	GIONEUT	2712
			R((119)=K(119)*Y(35)*Y(9)	GIONEUT	2713
			R((120)=K(120)*Y(36)	GIONEUT	2714
			R((121)=K(121)*Y(36)*Y(7)	GIONEUT	2715
			R((122)=K(122)*Y(56)*Y(18)	GIONEUT	2716
			R((123)=K(123)*Y(36)	GIONEUT	2717
			R((124)=K(124)*Y(31)*Y(9)	GIONEUT	2718
			R((125)=K(125)*Y(29)*Y(23)	GIONEUT	2719
			R((126)=K(126)*Y(29)*Y(8)	GIONEUT	2720
			R((127)=K(127)*Y(3)*Y(25)	GIONEUT	2721
			R((128)=K(128)*Y(7)*Y(7)*Y(57)	GIONEUT	2722
			R((129)=K(129)*Y(7)*Y(8)*Y(57)	GIONEUT	2723
			R((130)=K(130)*Y(7)*Y(9)	GIONEUT	2724
			R((131)=K(131)*Y(7)*Y(15)	GIONEUT	2725
			R((132)=K(132)*Y(7)*Y(14)	GIONEUT	2726
			R((133)=K(133)*Y(7)*Y(14)	GIONEUT	2727
			R((134)=K(134)*Y(11)*Y(14)	GIONEUT	2728
			R((135)=K(135)*Y(11)*Y(14)	GIONEUT	2729
			R((136)=K(136)*Y(11)*Y(8)*Y(57)	GIONEUT	2730
			R((137)=K(137)*Y(16)*Y(8)	GIONEUT	2731
			R((138)=K(138)*Y(16)*Y(18)	GIONEUT	2732
			R((139)=K(139)*Y(16)*Y(7)*Y(57)	GIONEUT	2733
			R((140)=K(140)*Y(10)*Y(9)	GIONEUT	2734
			R((141)=K(141)*Y(10)*Y(14)	GIONEUT	2735
			R((142)=K(142)*Y(18)*Y(9)	GIONEUT	2736
			R((143)=K(143)*Y(18)*Y(7)*Y(57)	GIONEUT	2737
			R((144)=K(144)*Y(12)*Y(12)	GIONEUT	2738
			R((145)=K(145)*Y(12)*Y(9)	GIONEUT	2739
			R((146)=K(146)*Y(19)*Y(11)	GIONEUT	2740
			R((147)=K(147)*Y(26)*Y(10)	GIONEUT	2741
			R((148)=K(148)*Y(26)*Y(7)*Y(57)	GIONEUT	2742
			R((149)=K(149)*Y(17)*Y(30)	GIONEUT	2743
			R((150)=K(150)*Y(5)*Y(30)	GIONEUT	2744
			R((151)=K(151)*Y(5)*Y(30)	GIONEUT	2745
			R((152)=K(152)*Y(4)*Y(30)	GIONEUT	2746
			R((153)=K(153)*Y(3)*Y(30)	GIONEUT	2747
			R((154)=K(154)*Y(6)*Y(30)	GIONEUT	2748
			R((155)=K(155)*Y(1)*Y(30)	GIONEUT	2749
			R((156)=K(156)*Y(40)*Y(30)	GIONEUT	2750
			R((157)=K(157)*Y(51)*Y(30)	GIONEUT	2751
			R((158)=K(158)*Y(52)*Y(30)	GIONEUT	2752
			R((159)=K(159)*Y(53)*Y(30)	GIONEUT	2753
			R((160)=K(160)*Y(8)*Y(30)	GIONEUT	2754
			R((161)=K(161)*Y(9)*Y(30)	GIONEUT	2755
			R((162)=K(162)*Y(4)*Y(8)*Y(57)	GIONEUT	2756
			R((163)=K(163)*Y(39)*Y(57)	GIONEUT	2757
			R((164)=K(164)*Y(39)*Y(27)	GIONEUT	2758
			R((165)=K(165)*Y(5)*Y(25)*Y(23)	GIONEUT	2759
			R((166)=K(166)*Y(5)*Y(13)*Y(57)	GIONEUT	2760
			R((167)=K(167)*Y(28)*Y(8)	GIONEUT	2761
			R((168)=K(168)*Y(3)*Y(23)	GIONEUT	2762
			R((169)=K(169)*Y(3)*Y(8)	GIONEUT	2763
			R((170)=K(170)*Y(41)*Y(13)*Y(57)	GIONEUT	2764
			R((171)=K(171)*Y(41)*Y(23)*Y(23)	GIONEUT	2765

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R(172)=K(172)*Y(42)*Y(13)*Y(57)          GIONEUT 2766
R(173)=K(173)*Y(43)*Y(13)*Y(57)          GIONEUT 2767
R(174)=K(174)*Y(43)*Y(57)                  GIONEUT 2768
R(175)=K(175)*Y(44)*Y(57)                  GIONEUT 2769
R(176)=K(176)*Y(45)*Y(57)                  GIONEUT 2770
R(177)=K(177)*Y(51)*Y(13)*Y(57)          GIONEUT 2771
R(178)=K(178)*Y(52)*Y(13)*Y(57)          GIONEUT 2772
R(179)=K(179)*Y(31)*Y(8)*Y(8)            GIONEUT 2773
R(180)=K(180)*Y(32)*Y(8)*Y(8)            GIONEUT 2774
R(181)=K(181)*Y(32)*Y(8)*Y(25)          GIONEUT 2775
R(182)=K(182)*Y(33)*Y(25)                GIONEUT 2776
R(183)=K(183)*Y(34)*Y(57)                GIONEUT 2777
R(184)=K(184)*Y(37)*Y(18)                GIONEUT 2778
R(185)=K(185)*Y(7)                      GIONEUT 2779
R(186)=K(186)*Y(8)                      GIONEUT 2780
R(187)=K(187)*Y(23)                      GIONEUT 2781
R(188)=K(188)*Y(18)                      GIONEUT 2782
R(189)=K(189)*Y(16)                      GIONEUT 2783
R(190)=K(190)*Y(14)                      GIONEUT 2784
R(191)=K(191)*Y(21)                      GIONEUT 2785
R(192)=K(192)*Y(8)                      GIONEUT 2786
R(193)=K(193)*Y(8)                      GIONEUT 2787
R(194)=K(194)*Y(9)                      GIONEUT 2788
R(195)=K(195)*Y(9)                      GIONEUT 2789
R(196)=K(196)*Y(13)                      GIONEUT 2790
R(197)=K(197)*Y(14)                      GIONEUT 2791
R(198)=K(198)*Y(12)                      GIONEUT 2792
R(199)=K(199)*Y(18)                      GIONEUT 2793
R(200)=K(200)*Y(19)                      GIONEUT 2794
R(201)=K(201)*Y(24)                      GIONEUT 2795
R(202)=K(202)*Y(55)                      GIONEUT 2796
R(203)=K(203)*Y(25)                      GIONEUT 2797
R(204)=K(204)*Y(25)                      GIONEUT 2798
R(205)=K(205)*Y(56)                      GIONEUT 2799
R(206)=K(206)*Y(7)                       GIONEUT 2800
R(207)=K(207)*Y(20)                      GIONEUT 2801
R(208)=K(208)*Y(23)                      GIONEUT 2802
R(209)=K(209)*Y(8)                       GIONEUT 2803
R(210)=K(210)*Y(8)                       GIONEUT 2804
R(211)=K(211)*Y(7)                       GIONEUT 2805
R(212)=K(212)*Y(7)                       GIONEUT 2806
R(213)=K(213)*Y(23)                      GIONEUT 2807
R(214)=K(214)*Y(8)                       GIONEUT 2808
R(215)=K(215)*Y(23)                      GIONEUT 2809
N=NREAC+NREAC2                           GIONEUT 2810
WRITE(6,20) (J,(CMB(I,J),I=1,6),K(J),R(J),J=1,MN) GIONEUT 2811
FORMAT(14,5A10,A8,9X,1PE12.5,9X,E12.5)      GIONEUT 2812
RETURN                                     GIONEUT 2813
END                                         GIONEUT 2814

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20 240 255 260 275

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1          SUBROUTINE HOUR(TIME,OUTP)
C THIS SUBROUTINE CONVERTS THE VALUE OF TIME IN SECONDS TO
C HOURS.
C
5          DIMENSION W(10),OUTP(1)
      DATA W/1H1,1H2,1H3,1H4,1H5,1H6,1H7,1H8,1H9,1H0/
      T=TIME+4.32E+04
      R=AMOD(T,8.64E+04)
      HR=R/3600.+0.00001
      IHR=HR
      FHR=HR-FLOAT(IHR)
      IFHR=FHR*60.01
      JL=4
10         DO 5 L=1,2
      IF(L.EQ.1) IW=IFHR
      IF(L.EQ.2) IW=IHR
      DO 5 M=1,2
      II=MOD(IW,10)
      IF(II.EQ.0) II=10
      OUTP(JL)=W(II)
      JL=JL-1
      IW=IW/10
      CONTINUE
20         RETURN
25         END
      GIONEUT 2815
      GIONEUT 2816
      GIONEUT 2817
      GIONEUT 2818
      GIONEUT 2819
      GIONEUT 2820
      GIONEUT 2821
      GIONEUT 2822
      GIONEUT 2823
      GIONEUT 2824
      GIONEUT 2825
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      GIONEUT 2829
      GIONEUT 2830
      GIONEUT 2831
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      GIONEUT 2833
      GIONEUT 2834
      GIONEUT 2835
      GIONEUT 2836
      GIONEUT 2837
      GIONEUT 2838
      GIONEUT 2839
      GIONEUT 2840

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